Final Fourth Quarter 2014 - Quarterly Groundwater Monitoring Report Inside Tunnel Wells

Red Hill Bulk Fuel Storage Facility
Joint Base Pearl Harbor-Hickam, Oahu, Hawaii

DOH Facility ID: 9-102271

DOH Release ID: 990051, 010011, 020028, and 140010

January 2015

Department of the Navy Naval Facilities Engineering Command, Hawaii 400 Marshall Road JBPHH HI 96860-3139



Contract Number N62742-12-D-1853, CTO 0002



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Prepared for:



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Prepared under:

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FINAL FOURTH QUARTER 2014 - QUARTERLY GROUNDWATER MONITORING REPORT INSIDE TUNNEL WELLS RED HILL BULK FUEL STORAGE FACILITY

Long-Term Groundwater and Soil Vapor Monitoring Red Hill Bulk Fuel Storage Facility Joint Base Pearl Harbor-Hickam, Oahu, Hawaii

Prepared for:

Department of the Navy
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Contract Number: N62742-12-D-1853 Contract Task Order: 0002

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Contract No. N62742-12-D-1853	Contract Task Order 0002
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ACRONYMS AND ABBREVIATIONS

ACRONYMS/

ABBREVIATIONS DEFINITION/MEANING

% percent

bgs below ground surface

COPC Contaminant of Potential Concern

DLNR State of Hawaii Department of Land and Natural Resources

DoD Department of Defense

DOH State of Hawaii Department of Health

DON Department of the Navy
EAL Environmental Action Level
EPA Environmental Protection Agency

ESI Environmental Science International, Inc.

F-76 Marine Diesel Fuel

ID Identification

JBPHH Joint Base Pearl Harbor-Hickam

JP-5 Jet Fuel Propellant-5
JP-8 Jet Fuel Propellant-8
LCS Laboratory Control Sample

LCSD Laboratory Control Sample Duplicate

LOD Limit of Detection
LOQ Limit of Quantitation
µg/L micrograms per liter

MS Matrix Spike

MSD Matrix Spike Duplicate

NAVFAC Naval Facilities Engineering Command

NAVSUP FLC Naval Supply Systems Command Fleet Logistics Center

PAH Polycyclic Aromatic Hydrocarbons

PARCCS Precision, Accuracy, Representativeness, Completeness, Comparability,

and Sensitivity

pH hydrogen activity

RHSF Red Hill Bulk Fuel Storage Facility

RPD Relative Percent Difference
SAP Sampling and Analysis Plan
SSRBL Site-Specific Risk-Based Level
TEC The Environmental Company, Inc.
TPH-d Total Petroleum Hydrocarbons as diesel
TPH-g Total Petroleum Hydrocarbons as gasoline

U.S. United States of America
UST Underground Storage Tank
VOC Volatile Organic Compound

WP Work Plan

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EXECUTIVE SUMMARY

This quarterly groundwater monitoring report presents the results of the fourth quarter 2014 groundwater sampling event, conducted on October 27 and 28, 2014, at the Red Hill Bulk Fuel Storage Facility [RHSF], Joint Base Pearl Harbor-Hickam [JBPHH], Hawaii. The RHSF is located in Halawa Heights on the Island of Oahu. There are 18 active and 2 inactive Underground Storage Tanks [USTs] located at the RHSF. The State of Hawaii Department of Health [DOH] Facility Identification [ID] number is 9-102271. The DOH Release ID numbers are 990051, 010011, 020028, and 140010.

The groundwater sampling was conducted as part of the long-term groundwater and soil vapor monitoring program at the RHSF and concurrent with release response activities initiated at Tank 5 in January 2014, under Naval Facilities Engineering Command [NAVFAC] Contract Number N62742-12-D-1853. The sampling was conducted in accordance with the approved 2012 Work Plan [WP]/Sampling and Analysis Plan [SAP] prepared by Environmental Science International, Inc. [ESI].

On October 27 and 28, 2014, ESI personnel collected groundwater samples from four monitoring wells at the RHSF (wells RHMW01, RHMW02, RHMW03, and RHMW05) and one sampling point at Red Hill Shaft (RHMW2254-01). A primary and duplicate sample were collected from well RHMW02. A summary of the analytical results is provided below.

- RHMW01 Total Petroleum Hydrocarbons as diesel fuel [TPH-d] (120 micrograms per liter [μg/L]) and lead (0.0976 μg/L) were the only analytes detected. The TPH-d concentration exceeded the DOH Environmental Action Level [EAL] for gross contamination, but was below the site-specific risk-based level [SSRBL] of 4,500 μg/L for Total Petroleum Hydrocarbons [TPH]. TPH-d concentrations in this well have shown an overall decreasing trend from a high of 1,500 μg/L in February 2005.
- RHMW02 TPH-d (2,000 μg/L in both primary and duplicate samples), Total Petroleum Hydrocarbons as gasoline [TPH-g] (57 and 53 μg/L), xylenes (0.32 and 0.29 μg/L), acenaphthene (0.53 μg/L in both primary and duplicate samples), ethylbenzene (0.15 and 0.14 μg/L), 1-methylnaphthalene (59 and 54 μg/L), 2-methylnaphthalene (43 and 36 μg/L), naphthalene (140 and 130 μg/L), and lead (0.165 μg/L in only the duplicate sample) were detected. TPH-d was detected at concentrations above the DOH EALs for both drinking water toxicity and gross contamination, but below half the SSRBL. The polycyclic aromatic hydrocarbons [PAHs] 1-methylnaphthalene, 2-methylnaphthalene, and naphthalene were detected at concentrations above the DOH EALs for both drinking water toxicity and gross contamination. The concentrations of 1-methylnaphthalene, 2-methylnaphthalene, and naphthalene increased to their highest levels in over 5 years. The concentrations of these PAHs are also significantly higher than they were in January 2014, when a high TPH-d spike was observed in this well.
- RHMW03 TPH-d (80 μg/L) was the only analyte detected. The concentration did not exceed the DOH EALs or the SSRBL.

- RHMW05 TPH-d (16 μ g/L) was the only analyte detected. The concentration did not exceed the DOH EALs or the SSRBL.
- RHMW2254-01 TPH-d (22 μg/L) and total lead (0.211 μg/L) were the only analytes detected. The concentrations did not exceed their respective DOH EAL.

During this quarterly event, the TPH-d concentration in RHMW01 exceeded the DOH EAL for gross contamination and the concentrations of TPH-d, 1-methylnaphthalene, 2-methylnaphthalene, and naphthalene in RHMW02 exceeded the DOH EALs for both gross contamination and drinking water toxicity. Groundwater contaminant concentrations in wells RHMW03, RHMW05, and RHMW2254-01 remained at low concentrations and did not change significantly from the previous event, or were not detected.

Concentrations of 1-methylnaphthalene, 2-methylnaphthalene, and naphthalene in RHMW02 have shown an increasing trend since March 5, 2014, but have remained below the historic maximums detected in the well. These PAH concentrations were an order of magnitude higher than they were during January 2014, when a TPH-d spike was observed but PAH concentrations remained below DOH EALs for gross contamination. Based on the groundwater monitoring data, TPH-d and PAH concentrations in RHMW02 appear to be decoupled. All other analytical results were generally consistent with historical data.

Based on the groundwater monitoring results and the reported release at Tank 5 in January 2014, continued groundwater monitoring at the wells inside the RHSF tunnel is recommended. The next quarterly event is tentatively scheduled for January 2015.

SECTION 1 – INTRODUCTION

This quarterly groundwater monitoring report presents the results of the fourth quarter 2014 groundwater sampling conducted on October 27 and 28, 2014, at the RHSF, JBPHH. The RHSF is located in Halawa Heights on the Island of Oahu. The purpose of the sampling is to (1) assess the condition of groundwater beneath the RHSF with respect to chemical constituents associated with jet fuel propellant and marine diesel fuel, and (2) to ensure the Navy remains in compliance with DOH UST release response requirements as described in Hawaii Administrative Rules Chapter 11-281 Subchapter 7, *Release Response Action* (DOH, 2013). The DOH Facility ID number for the RHSF is 9-102271. The DOH Release ID numbers are 990051, 010011, 020028, and 140010.

The groundwater sampling was conducted as part of the long-term groundwater and soil vapor monitoring program at the RHSF, under NAVFAC Contract Number N62742-12-D-1853. The sampling was conducted in accordance with the approved WP/SAP prepared by ESI (ESI, 2012).

1.1 SITE DESCRIPTION

The RHSF is located on federal government land (zoned F1- Military and Federal), located in Halawa Heights, approximately 2.5 miles northeast of Pearl Harbor (Figure 1). It is located on a low ridge on the western edge of the Koolau Mountain Range that divides Halawa Valley from Moanalua Valley. The RHSF is bordered on the north by Halawa Correctional Facility and private businesses, on the southwest by the United States of America [U.S.] Coast Guard reservation, on the south by residential neighborhoods, and on the east by Moanalua Valley. A quarry is located less than a quarter mile away to the northwest. The RHSF occupies 144 acres of land and the majority of the site is at an elevation of approximately 200 to 500 feet above mean sea level.

The RHSF contains 18 active and 2 inactive USTs that are operated by Naval Supply Systems Command Fleet Logistics Center [NAVSUP FLC] Pearl Harbor (formerly Fleet and Industrial Supply Center). Each UST has a capacity of approximately 12.5 million gallons. The RHSF is located approximately 100 feet above the basal aquifer. The USTs contain Jet Fuel Propellant-5 [JP-5], Jet Fuel Propellant-8 [JP-8], and Marine Diesel Fuel [F-76]. The current status of the USTs are summarized in Table 1.1.

Four groundwater monitoring wells (wells RHMW01, RHMW02, RHMW03, and RHMW05) and one sampling point at Red Hill Shaft (RHMW2254-01) are located within the RHSF lower access tunnel (Figure 2). Three groundwater monitoring wells (wells HDMW2253-03, OWDFMW01, and RHMW04) are located outside of the RHSF tunnel system. Monitoring data for the three wells located outside the tunnel are included in a separate report.

As noted, monitoring wells RHMW01, RHMW02, RHMW03, and RHMW05 are located inside the underground tunnels. Sampling point RHMW2254-01 is located inside the infiltration gallery

of the Department of the Navy [DON] drinking water supply Well 2254-01. The DON Well 2254-01 is located approximately 2,400 feet downgradient of the USTs and provides potable water to the JBPHH Water System, which serves approximately 65,200 military customers. NAVFAC Public Works Department operates the infiltration gallery and DON Well 2254-01.

TABLE 1.1
Current Status of the USTs
Red Hill Bulk Fuel Storage Facility
October 2014 Quarterly Monitoring Report

Tank Identification	Fuel Type	Status	Capacity
F-1	None	Inactive	12.5 million gallons
F-2	JP-8	Active	12.5 million gallons
F-3	JP-8	Active	12.5 million gallons
F-4	JP-8	Active	12.5 million gallons
F-5	JP-8	Active	12.5 million gallons
F-6	JP-8	Active	12.5 million gallons
F-7	JP-5	Active	12.5 million gallons
F-8	JP-5	Active	12.5 million gallons
F-9	JP-5	Active	12.5 million gallons
F-10	JP-5	Active	12.5 million gallons
F-11	JP-5	Active	12.5 million gallons
F-12	JP-5	Active	12.5 million gallons
F-13	F-76	Active	12.5 million gallons
F-14	F-76	Active	12.5 million gallons
F-15	F-76	Active	12.5 million gallons
F-16	F-76	Active	12.5 million gallons
F-17	JP-5	Active	12.5 million gallons
F-18	JP-5	Active	12.5 million gallons
F-19	None	Inactive	12.5 million gallons
F-20	JP-5	Active	12.5 million gallons

F-76 Marine Diesel Fuel

JP-5 Jet Fuel Propellant-5

JP-8 Jet Fuel Propellant-8

1.2 PHYSICAL SETTING

Climatological conditions in the area of the RHSF consist of warm to moderate temperatures and low to moderate rainfall. The RHSF is leeward of the prevailing northeasterly trade winds. The average annual precipitation is approximately 40 inches, which occurs mainly between November and April (State of Hawaii Department of Land and Natural Resources [DLNR], 1986). Annual pan evaporation is approximately 75 inches (DLNR, 1985). Average temperatures range from the low 60's to high 80's (degrees Fahrenheit) (Atlas of Hawaii, 1983).

Oahu consists of the eroded remnants of two shield volcanoes, Waianae and Koolau. The RHSF is located on the southwest flank of the Koolau volcanic shield. Lavas erupted during the shield-building phase of the volcano belong to the *Koolau Volcanic Series* (Stearns and Vaksvik, 1935). Following formation of the Koolau shield, a long period of volcanic quiescence occurred, during which the shield was deeply eroded. Following this erosional period, eruptive activity resumed. Lavas and pyroclastic material erupted during this period belong to the *Honolulu*

1-2

Volcanic Series (Stearns and Vaksvik, 1935).

In the immediate area of the RHSF, Koolau Volcanic Series lavas dominate, although there are consolidated and unconsolidated non-calcareous deposits in the vicinity that consist of alluvium generated during erosion of the Koolau volcanic shield. South-southwest of the RHSF, and in isolated exposures to the west, are pyroclastic deposits formed during eruptions from three Honolulu Volcanic Series vents, Salt Lake, Aliamanu, and Makalapa (Stearns and Vaksvik, 1935). Based on established geology and records of wells drilled at the RHSF (Stearns and Vaksvik, 1938), the RHSF is underlain by Koolau Volcanic Series basalts. The area of the RHSF is classified as *Rock Land*, where 25-90% of the land surface is covered by exposed rock and there are only shallow soils (Foote, et al., 1972).

Groundwater in Hawaii exists in two principal aquifer types. The first and most important type, in terms of drinking water resources, is the basal aquifer. The basal aquifer exists as a lens of fresh water floating on and displacing seawater within the pore spaces, fractures, and voids of the basalt that forms the underlying mass of each Hawaiian island. In parts of Oahu, groundwater in the basal aquifer is confined by the overlying caprock and is under pressure. Waters that flow freely to the surface from wells that tap the basal aquifer are referred to as artesian.

The second type of aquifer is the caprock aquifer, which consists of various kinds of unconfined and semi-confined groundwater. Commonly, the caprock consists of a thick sequence of nearly impermeable clays, coral, and basalt, that separates the caprock aquifer from the basal aquifer. The impermeable nature of these materials and the artesian nature of the basal aquifer severely restrict the downward migration of groundwater from the upper caprock aquifer. However, in the area of the RHSF, there is no discernible caprock.

Groundwater in the area of the RHSF is part of the *Waimalu Aquifer System* of the *Pearl Harbor Aquifer Sector*. The aquifer is classified as a basal, unconfined, flank-type; and is currently used as a drinking water source. The aquifer is considered fresh, with less than 250 milligrams per liter of chloride, and is considered an irreplaceable resource with a high vulnerability to contamination (Mink and Lau, 1990).

The nearest drinking water supply well is DON Well 2254-01, located in the infiltration gallery within the RHSF lower tunnel. The DON Well 2254-01 is located approximately 2,400 feet hydraulically and topographically downgradient of the USTs (Figure 2).

1.3 BACKGROUND

The RHSF was constructed by the U.S. Government in the early 1940s. Twenty USTs and a series of tunnels were constructed. The USTs were constructed of steel, and in the past have stored DON special fuel oil, DON distillate, aviation gasoline, and motor gasoline (Environet, 2010). The tanks currently contain JP-5, JP-8, and F-76. The fueling system is a self-contained underground unit that was installed into native rock comprised primarily of basalt with some

interbedded tuffs and breccias (Environet, 2010). Each UST measures approximately 250 feet in height and 100 feet in diameter. The upper domes of the tanks lie at a depth varying between 100 feet and 200 feet below ground surface [bgs].

In 1998, Earth Tech conducted a Phase II Remedial Investigation/Feasibility Study for the Oily Waste Disposal Facility located within the RHSF. The study included the installation of well OWDFMW01 (which was originally MW08) (Earth Tech, 1999).

In February 2001, the DON installed groundwater monitoring well RHMW01 to monitor for contamination in the basal aquifer beneath the RHSF. Well RHMW01 was installed approximately 100 feet below grade within the lower access tunnel. The depth to water was measured at 86 feet below the tunnel floor at the time of the well completion. In February 2001, a groundwater sample was collected from the well. TPH and total lead were detected in the sample. Total lead was detected at a concentration above the DOH Tier 1 groundwater action level of 5.6 µg/L (The Environmental Company, Inc. [TEC], 2009; DOH, 2000).

In 2005, the RHSF groundwater monitoring program was initiated. It involved routine groundwater sampling of well RHMW01 and sampling point RHMW2254-01. Samples were collected in February, June, September, and December of 2005. Lead was detected at concentrations above the DOH Tier 1 action level of 5.6 µg/L in samples collected in February and June. The samples collected in February and June were not filtered prior to analysis, whereas the samples collected in September and December were filtered prior to analysis. Since the samples collected in February and June were not filtered prior to analysis, the lead results were not considered appropriate for a risk assessment (TEC, 2008).

Between June and September 2005, TEC installed three additional groundwater monitoring wells (wells RHMW02, RHMW03, and RHMW04) (TEC, 2008). Well RHMW04 was installed hydraulically upgradient of the USTs to provide background geochemistry information for water moving through the basal aquifer beneath the RHSF. Wells RHMW02 and RHMW03 were installed approximately 125 feet below grade within the RHSF lower tunnel and well RHMW04 was installed to a depth of approximately 300 feet bgs outside of the RHSF tunnels. In September 2005, groundwater samples were collected from the three newly installed groundwater monitoring wells (wells RHMW02, RHMW03, and RHMW04) along with the existing well RHMW01 and sampling point RHMW2254-01. The contaminants of potential concern [COPCs] with concentrations exceeding current DOH EALs are summarized below.

- RHMW01 TPH-d was detected at concentrations above the DOH EALs.
- **RHMW02** TPH-g, TPH-d, naphthalene, trichloroethylene, 1-methylnaphthalene, and 2-methylnaphthalene were detected at concentrations above their respective DOH EAL.
- RHMW03 TPH-d was detected at concentrations above the DOH EALs.

In 2006, TEC installed dedicated sampling pumps in the four wells (wells RHWM01, RHWM02, RHMW03, and RHWM04) and one sampling point (RHMW2254-01). In July and December of

2006, groundwater samples were collected from the four wells and the sampling point. The COPCs with concentrations exceeding current DOH EALs are summarized below.

- RHMW01 TPH-d and naphthalene were detected at concentrations above their respective DOH EAL.
- RHMW02 TPH-g, TPH-d, and naphthalene were detected at concentrations above their respective DOH EAL.
- RHMW03 TPH-d was detected at concentrations above the DOH EALs.

In 2007, SSRBLs were established for TPH (4,500 μ g/L) and benzene (750 μ g/L) based on the solubility in water of JP-5 and JP-8 (TEC, 2007). Groundwater samples were collected from wells RHWM01, RHWM02, and RHMW03, and sampling point RHMW2254-01. Samples were collected in March, June, and September of 2007. The COPCs with concentrations exceeding current DOH EALs are summarized below.

- RHMW01 TPH-d was detected at concentrations above the DOH EALs, but below the SSRBL.
- RHMW02 TPH-g, TPH-d, naphthalene, 1-methylnaphthalene, and 2-methylnaphthalene were detected at concentrations above their respective DOH EAL. The TPH-d concentrations were below the SSRBL.
- RHMW03 TPH-d was detected at concentrations above the DOH EALs, but below the SSRBL.

In 2008, groundwater samples were collected from wells RHWM01, RHWM02, and RHMW03, and sampling point RHMW2254-01. Samples were collected in January, April, July, and October of 2008. The COPCs with concentrations exceeding current DOH EALs are summarized below. In addition, a groundwater protection plan (TEC, 2008) was prepared.

- RHMW01 TPH-d was detected at concentrations above the DOH EALs, but below the SSRBL.
- RHMW02 TPH-d, naphthalene, 1-methylnaphthalene, and 2-methylnaphthalene were detected at concentrations above their respective DOH EAL. The TPH-d concentrations detected in October 2008 were also above the SSRBL.
- RHMW03 TPH-d was detected at concentrations above the DOH EALs, but below the SSRBL.

In April 2009, groundwater monitoring well RHMW05 was installed downgradient of the USTs, within the lower access tunnel between RHMW01 and RHMW2254-01. It was installed to identify the extent of contamination hydraulically downgradient of the USTs. Well RHMW05 was added to the quarterly groundwater sampling program. In 2009, quarterly groundwater samples were collected from wells RHWM01, RHWM02, RHMW03, and RHMW05, and sampling point RHMW2254-01. Samples were collected in February, May, July, and October of 2009. The

COPCs with concentrations exceeding current DOH EALs are summarized below. In addition, the Groundwater Protection Plan was revised to include well RHMW05.

- RHMW01 TPH-d and 1-methylnaphthalene were detected at concentrations above their respective DOH EAL. The TPH-d concentrations were below the SSRBL.
- RHMW02 TPH-d, naphthalene, 1-methylnaphthalene, and 2-methylnaphthalene were detected at concentrations above their respective DOH EAL. The TPH-d concentrations were below the SSRBL.
- RHMW03 TPH-d was detected at a concentration above the DOH EALs, but below the SSRBL.
- RHMW05 TPH-d was detected at a concentration above the DOH EALs, but below the SSRBL.

In 2010, groundwater samples were collected from wells RHWM01, RHWM02, RHMW03, and RHMW05, and sampling point RHMW2254-01. Samples were collected in January, April, July, and October. The COPCs with concentrations exceeding current DOH EALs are summarized below.

- RHMW01 TPH-d was detected at concentrations above the DOH EALs, but below the SSRBL.
- RHMW02 TPH-g, TPH-d, naphthalene, and 1-methylnaphthalene were detected at concentrations above their respective DOH EAL. The TPH-d concentrations were below the SSRBL.
- RHMW03 TPH-d was detected at a concentration above the DOH EALs, but below the SSRBL.
- RHMW05 TPH-d was detected at a concentration above the DOH EALs, but below the SSRBL.

In 2011, quarterly groundwater samples were collected from wells RHWM01, RHWM02, RHMW03, and RHMW05, and sampling point RHMW2254-01. Samples were collected in January, April, July, and October. In a Fall 2011 update, the DOH EALs were revised. The drinking water toxicity EAL for TPH-d decreased from 210 to 190 μ g/L (DOH, 2011). The COPCs with concentrations exceeding current DOH EALs are summarized below.

- RHMW01 TPH-d was detected at concentrations above the DOH EALs, but below the SSRBL.
- RHMW02 TPH-d, naphthalene, indeno[1,2,3-cd]pyrene, and 1-methylnaphthalene were
 detected at concentrations above their respective DOH EAL. The TPH-d concentrations
 were below the SSRBL.

In 2012, quarterly groundwater samples were collected from wells RHWM01, RHWM02, RHMW03, and RHMW05, and sampling point RHMW2254-01. Samples were collected in

February, April, July, and November. The COPCs with concentrations exceeding current DOH EALs are summarized below.

- RHMW01 TPH-d was detected at concentrations above the DOH EALs, but below the SSRBL.
- RHMW02 TPH-d, TPH-g, naphthalene, 1-methylnaphthalene, and 2-methylnaphthalene were detected at concentrations above their respective DOH EAL. The TPH-d concentrations were below the SSRBL.

In 2013, quarterly groundwater samples were collected from wells RHWM01, RHWM02, RHMW03, and RHMW05, and sampling point RHMW2254-01. Samples were collected in January, April, July, and October. The COPCs with concentrations exceeding current DOH EALs are summarized below.

- RHMW01 TPH-d was detected at concentrations above the DOH EALs, but below the SSRBL.
- RHMW02 TPH-d, TPH-g, naphthalene, 1-methylnaphthalene, and 2-methylnaphthalene were detected at concentrations above their respective DOH EAL. The TPH-d concentrations were below the SSRBL.

In January 2014, quarterly groundwater samples were collected from wells RHMW01, RHMW02, RHMW03, RHMW05, and sampling point RHMW2254-01. The COPCs with concentrations exceeding current DOH EALs are summarized below.

- RHMW01 TPH-d was detected at a concentration above the DOH EALs, but below the SSRBL.
- RHMW02 TPH-d was detected at concentrations above both DOH EALs, and naphthalene
 and 1-methylnaphthalene were detected at concentrations above only the DOH EAL for
 drinking water toxicity. The TPH-d concentrations exceeded the SSRBL.

Between January and April 2014, additional groundwater sampling was conducted at wells RHMW01, RHMW02, RHMW05, and sampling point RHMW2254-01 in response to a reported release from Tank 5. The COPCs with concentrations exceeding current DOH EALs are summarized below.

- RHMW01 TPH-d was detected at concentrations above the DOH EALs, but below the SSRBL.
- RHMW02 TPH-d was detected at concentrations above both DOH EALs, and naphthalene
 and 1-methylnaphthalene were detected at concentrations above only the DOH EAL for
 drinking water toxicity. The TPH-d concentrations were below the SSRBL.

In April 2014, quarterly groundwater samples were collected from wells RHMW01, RHMW02, RHMW03, RHMW05, and sampling point RHMW2254-01. The COPCs with concentrations exceeding current DOH EALs are summarized below.

 RHMW02 – TPH-d and naphthalene were detected at concentrations above both DOH EALs. 1-Methylnaphthalene was detected at concentrations above only the DOH EAL for drinking water toxicity. The TPH-d concentrations were below the SSRBL.

In May and June 2014, additional groundwater sampling was conducted at wells RHMW01, RHMW02, RHMW05, and sampling point RHMW2254-01 in response to a reported release from Tank 5. The COPCs with concentrations exceeding current DOH EALs are summarized below.

 RHMW02 – TPH-d, naphthalene, and 1-methylnaphthalene were detected at concentrations above their respective DOH EAL. The TPH-d concentrations were below the SSRBL.

In July 2014, quarterly groundwater samples were collected from wells RHMW01, RHMW02, RHMW03, RHMW05, and sampling point RHMW2254-01. The COPCs with concentrations exceeding current DOH EALs are summarized below.

 RHMW02 – TPH-d, 1-methylnaphthalene, and naphthalene were detected at concentrations above their respective DOH EAL. 2-Methylnaphthalene was detected at concentrations above only the DOH EAL for gross contamination. The TPH-d concentrations were below the SSRBL.

1.3.1 Previous Reports

The following groundwater monitoring reports were previously submitted to the DOH:

- 1. Groundwater Sampling Report, First Quarter 2005 (submitted April 2005).
- Groundwater Sampling Report, Second Quarter 2005 (submitted August 2005).
- 3. Groundwater Sampling Report, Third Quarter 2005 (submitted November 2005).
- 4. Groundwater Sampling Report, Fourth Quarter 2005 (submitted February 2006).
- Groundwater Monitoring Results, July 2006 (submitted September 2006).
- Groundwater Monitoring Results, December 2006 (submitted January 2007).
- 7. Groundwater Monitoring Results, March 2007 (submitted May 2007).
- 8. Groundwater Monitoring Results, June 2007 (submitted August 2007).
- 9. Groundwater Monitoring Results, September 2007 (submitted October 2007).
- 10. Groundwater Monitoring Report, January 2008 (submitted March 2008).
- 11. Groundwater Monitoring Report, April 2008 (submitted May 2008).

- 12. Groundwater Monitoring Report, July 2008 (submitted October 2008).
- 13. Groundwater Monitoring Report, October and December 2008 (submitted February 2009).
- 14. Groundwater Monitoring Report, February 2009 (submitted May 2009).
- 15. Groundwater Monitoring Report, May 2009 (submitted July 2009).
- 16. Groundwater Monitoring Report, July 2009 (submitted September 2009).
- 17. Groundwater Monitoring Report, October 2009 (submitted December 2009).
- 18. Groundwater Monitoring Report, January, February, and March 2010 (submitted April 2010).
- 19. Groundwater Monitoring Report, April 2010 (submitted May 2010).
- 20. Groundwater Monitoring Report, July 2010 (submitted August 2010).
- 21. Groundwater Monitoring Report, October 2010 (submitted December 2010).
- 22. Groundwater Monitoring Report, January 2011 (submitted March 2011).
- 23. Groundwater Monitoring Report, April 2011 (submitted June 2011).
- 24. Groundwater Monitoring Report, July 2011 (submitted September 2011).
- 25. Groundwater Monitoring Report, October 2011 (submitted December 2011).
- 26. Groundwater Monitoring Report, January-February 2012 (submitted March 2012).
- 27. Groundwater Monitoring Report, April 2012 (Submitted July 2012).
- 28. Groundwater Monitoring Report, October 2012 (Submitted January 2013).
- 29. Groundwater Monitoring Report, January 2013 (Submitted April 2013).
- 30. Groundwater Monitoring Report, April 2013 (Submitted July 2013).
- 31. Groundwater Monitoring Report, July 2013 (Submitted September 2013).
- 32. Groundwater Monitoring Report, October 2013 (Submitted January 2014).
- 33. Groundwater Sampling Report for Additional Sampling, January 2014 (submitted January 2014).
- 34. Groundwater Monitoring Report, January 2014 (Submitted April 2014).
- 35. Groundwater Sampling Report for Tank 5 Release Response on March 5 and 6, 2014 (submitted March 2014).
- 36. Groundwater Sampling Report for Tank 5 Release Response on March 10, 2014 (submitted March 2014).
- 37. Groundwater Sampling Report for Tank 5 Release Response on March 25 and 26, 2014 (submitted April 2014).

- 38. Groundwater Sampling Report for Tank 5 Release Response on April 7, 2014 (submitted April 2014).
- 39. Groundwater Monitoring Report, April 2014 (Submitted June 2014).
- 40. Groundwater Sampling Report for Tank 5 Release Response on May 27 and 28, 2014 (submitted June 2014).
- 41. Groundwater Sampling Report for Tank 5 Release Response on June 23 and 24, 2014 (submitted July 2014).
- 42. Groundwater Monitoring Report, July 2014 (Submitted September 2014).

SECTION 2 – GROUNDWATER SAMPLING

On October 27 and 28, 2014, ESI personnel collected groundwater samples from four monitoring wells at the RHSF (wells RHMW01, RHMW02, RHMW03, and RHMW05) and one sampling point at Red Hill Shaft (RHMW2254-01). The samples were collected in accordance with the 2012 WP/SAP. The WP/SAP is consistent with DOH UST release response requirements (DOH, 2000); DON Procedure I-C-3, *Monitoring Well Sampling* (DON, 2007); and the RHSF Groundwater Protection Plan (TEC, 2008). Prior to purging and sampling, the depth to groundwater and the depth to the bottoms of the wells were measured using a Geotech oil/water interface probe. No measurable product, sheen, or petroleum hydrocarbon odor was detected in any of the wells.

2.1 GROUNDWATER SAMPLING

Prior to collecting groundwater samples, the monitoring wells were purged of standing water in the well casings. Each well contains a dedicated bladder pump which was used to purge the well and to collect samples. To operate the pump, a portable air compressor with an in-line filter was connected to a QED MP50 MicroPurge® Basics Controller box, which was then connected to the pump. The compressor was turned on to power the pump and the controller was used to adjust the pumping rate to less than one liter of water per minute.

Water quality parameters were monitored on a periodic basis during well purging. The water quality parameters that were measured included hydrogen activity [pH], temperature, conductivity, dissolved oxygen, and oxidation reduction potential. The water quality parameters were evaluated to assess whether the natural characteristics of the aquifer formation water were present within the monitoring wells before collecting the samples. At least four readings were collected during the purging process. Purging was considered complete when at least three consecutive water quality measurements stabilized within approximately 10%. The readings were recorded on groundwater monitoring logs. The groundwater monitoring logs are included in Appendix A. In addition, field notes were taken to document the sampling event. The field notes are included in Appendix B.

When the water quality parameters stabilized, groundwater samples were collected from the wells using the bladder pumps. The groundwater samples were collected no more than two hours after purging was completed to decrease groundwater interaction with the monitoring well casing and atmosphere. Prior to collecting the sample, the water level in the monitoring wells was measured and recorded to ensure that excessive drawn down had not occurred. The groundwater samples were collected at flow rates of approximately 0.17 to 0.5 liters per minute. Samples collected for dissolved lead analysis were filtered in the field using new, 0.45-micron filters.

All samples were labeled and logged on the Sample Inventory Log, placed in Ziploc™ bags and sealed, custody sealed, sealed with tape, placed in a cooler with wet ice, and logged onto the Chain-of-Custody form. The samples were labeled and logged in accordance with Project

Procedures Manual Standard Operating Procedure III-E, Record Keeping, Sample Labeling, and Chain-of-Custody Procedures (DON, 2007). All samples were shipped under Chain-of-Custody to the analytical laboratory and analyzed for the COPCs as described in Section 2.2.

2.2 ANALYTICAL RESULTS

The samples were analyzed for TPH-d using U.S. Environmental Protection Agency [EPA] Method 8015M, TPH-g and Volatile Organic Compounds [VOCs] using EPA Method 8260B, PAHs using EPA Method 8270C SIM, dissolved lead using EPA Method 6020, and total lead using EPA Method 200.8. The sample collected from sampling point RHMW2254-01 was analyzed for total lead (unfiltered) as DON Well 2254-01 is a drinking water supply well. A copy of the laboratory report is included as Appendix C.

Analytical results were compared to the DOH EALs for drinking water toxicity and gross contamination. Analytical results for wells RHMW01, RHMW02, RHMW03, and RHMW05 were also compared to the SSRBLs for TPH (4,500 μ g/L) and benzene (750 μ g/L), established in the RHSF Final Groundwater Protection Plan (TEC, 2008). The results of the fourth quarter groundwater sampling event are summarized in Table 2.1 and described below.

- RHMW01 TPH-d (120 µg/L) and lead (0.0976 µg/L) were the only analytes detected. The
 TPH-d concentration exceeded the DOH EAL for gross contamination, but was below the
 SSRBL of 4,500 µg/L for TPH. None of the other chemical constituents analyzed for were
 detected at concentrations above their respective DOH EAL.
- RHMW02 TPH-d (2,000 μg/L), TPH-g (57 and 53 μg/L), xylenes (0.32 and 0.29 μg/L), acenaphthene (0.53 μg/L), ethylbenzene (0.15 and 0.14 μg/L), 1-methylnaphthalene (59 and 54 μg/L), 2-methylnaphthalene (43 and 36 μg/L), and naphthalene (140 and 130 μg/L) were detected in both the primary and duplicate samples collected. Lead (0.165 μg/L) was only detected in the duplicate sample. TPH-d, 1-methylnaphthalene, 2-methylnaphthalene, and naphthalene were detected at concentrations above their respective DOH EALs for both drinking water toxicity and gross contamination. However, the TPH-d concentrations did not exceed the SSRBL.
- RHMW03 TPH-d (80 μg/L) was the only analyte detected. The concentration did not exceed the DOH EAL or the SSRBL.
- RHMW05 TPH-d (16 μ g/L) was the only analyte detected. The concentration did not exceed the DOH EAL or the SSRBL.
- RHMW2254-01 TPH-d (22 μg/L) and total lead (0.211 μg/L) were the only analytes detected. The concentrations did not exceed their respective DOH EAL.

2.3 GROUNDWATER CONTAMINANT TRENDS

The historical groundwater contaminant concentrations for COPCs that exceeded the DOH EALs or SSRBLs are illustrated in Appendix D. A table of cumulative historical groundwater

results is included as Appendix E. A summary of groundwater contaminant trends is provided below.

- RHMW01 COPCs detected during this round of quarterly sampling are consistent with the
 historical data for RHMW01. TPH-d has historically been detected at concentrations above
 the DOH EAL for both drinking water toxicity and gross contamination. TPH-d
 concentrations have shown an overall decreasing trend from a high of 1,500 μg/L in
 February 2005. The TPH-d concentration exceeded the DOH EAL for gross contamination
 for the first time since April 7, 2014.
- RHMW02 COPCs detected during this round of quarterly sampling are consistent with the historical data for RHMW02. TPH-g, TPH-d, 1-methylnaphthalene, 2-methylnaphthalene, and naphthalene have historically been detected at concentrations above the DOH EALs. During the October 2014 event, concentrations of TPH-d, 1-methylnaphthalene, 2-methylnaphthalene, and naphthalene approximately doubled from the previous event in July 2014 and remained above DOH EALs. The concentrations of 1-methylnaphthalene, 2-methylnaphthalene, and naphthalene increased to their highest levels in over 5 years. The concentrations of these PAHs are also significantly higher than they were in January 2014, when a high TPH-d spike was observed in this well. The concentrations of TPH-g remained below the DOH EALs for gross contamination and drinking water toxicity and were comparable to the concentrations detected during the previous event. Trichloroethylene was detected once in RHMW02 in September 2005 in the primary sample at a concentration above the DOH EAL for drinking water toxicity; however, trichloroethylene was not detected in the duplicate sample, and this may have been an anomalous result.
- RHMW03 COPCs detected during this round of quarterly sampling are consistent with the historical data for RHMW03. TPH-d has historically been detected at concentrations above the DOH EALs; however, it has not been detected at concentrations above the DOH EALs since October 2010.
- RHMW05 COPCs detected during this round of quarterly sampling are consistent with the
 historical data for RHMW05. TPH-d has historically been detected in RHMW05 at
 concentrations above the DOH EAL for both drinking water toxicity and gross contamination;
 however, it has not been detected at concentrations above the DOH EALs since January
 2010.
- RHMW2254-01 COPCs detected during this round of quarterly sampling are consistent
 with the historical data for RHMW2254-01. Although the method reporting limits for TPH-d
 exceeded one or both DOH EALs for drinking water toxicity and gross contamination
 between May 2009 and July 2010, TPH-d was last detected in RHMW2254-01 at a
 concentration above the DOH EAL for gross contamination in January 2008.

2.4 WASTE DISPOSAL

The purged groundwater and decontamination water generated during sampling of the inside tunnel was placed in a 55-gallon drum along with the purged water and decontamination water from the outside tunnel wells. The drum is currently stored onsite at ADIT 3 on top of a

secondary containment spill pallet and covered by a tarp. There is a non-hazardous label affixed to the drum with all pertinent information relating to its generation. The drum will be used for future sampling events and will be properly disposed of once it has been filled.

Contract No. N62742-12-D-1853 Contract Task Order 0002

TABLE 2.1

Analytical Results for Groundwater Sampling (October 27 and 28, 2014) Red Hill Bulk Fuel Storage Facility October 2014 Quarterly Monitoring Report

		DOH E	RHMW2254-01 (ES117)				RHWM01 (ES113)					RHMW02 (ES114)						RHW	VM03 (ES	116)		RHMW05 (ES118)						
Method	Chemical	Drinking Water Toxicity	Gross Contamination	Results	Q	LOQ	LOD	DL	Results	Q	LOQ	LOD	DL	Results	Q	LOQ	LOD	DL	Results	Q	LOQ	LOD	DL	Results	Q	LOQ	LOD	DL
EPA 8015B	TPH-d	190	100	22	HD,J	25	12	11	120	HD,J	25	12	11	2,000	HD,J	25	12	11	80 N.D.	HD,J U	25	12	11	16 N.D.	HD,J	25	12	11
EPA 8260B	TPH-g Acenaphthene	100 370	100 20	N.D. N.D.	U	50 0.19	30 0.049	26 0.026	N.D. N.D.	U	50 0.21	30 0.052	26 0.028	57 0.53		50 0.19	30 0.047	26 0.025	N.D. N.D.	U	50 0.21	30 0.054	26 0.029	N.D. N.D.	U	50 0.19	30 0.048	26 0.026
	Acenaphthylene	240	2,000	N.D.	U	0.19	0.049	0.043	N.D.	U	0.21	0.052	0.026	N.D.	U	0.19	0.047	0.042	N.D.	U	0.21	0.054	0.048	N.D.	U	0.19	0.048	0.020
	Anthracene	1,800	22	N.D.	Ü	0.19	0.049	0.028	N.D.	Ü	0.21	0.052	0.030	N.D.	Ü	0.19	0.047	0.027	N.D.	Ü	0.21	0.054	0.031	N.D.	Ü	0.19	0.048	0.028
	Benzo[a]anthracene	0.092	4.7	N.D.	U	0.19	0.049	0.032	N.D.	U	0.21	0.052	0.034	N.D.	U	0.19	0.047	0.031	N.D.	U	0.21	0.054	0.035	N.D.	U	0.19	0.048	0.031
	Benzo[g,h,i]perylene	1,500	0.13	N.D.	U	0.19	0.097	0.080	N.D.	U	0.21	0.10	0.085	N.D.	U	0.19	0.095	0.078	N.D.	U	0.21	0.11	0.088	N.D.	U	0.19	0.096	0.079
	Benzo[a]pyrene	0.2	0.81	N.D.	U	0.19	0.049	0.022	N.D.	U	0.21	0.052	0.023	N.D.	U	0.19	0.047	0.021	N.D.	U	0.21	0.054	0.024	N.D.	U	0.19	0.048	0.021
	Benzo[b]fluoranthene	0.092	0.75	N.D.	U	0.19	0.049	0.017	N.D.	U	0.21	0.052	0.018	N.D.	U	0.19	0.047	0.017	N.D.	U	0.21	0.054	0.019	N.D.	U	0.19	0.048	0.017
	Benzo[k]fluoranthene Chrysene	0.92 9.2	0.4	N.D. N.D.	U	0.19 0.19	0.049 0.049	0.030 0.024	N.D. N.D.	U	0.21 0.21	0.052 0.052	0.032 0.026	N.D. N.D.	U	0.19 0.19	0.047 0.047	0.029 0.024	N.D. N.D.	U	0.21 0.21	0.054 0.054	0.033 0.027	N.D. N.D.	U	0.19 0.19	0.048	0.030 0.024
EPA 8270C	Dibenzo[a,h]anthracene ¹	0.0092	0.52	N.D.	U	0.19	0.049	0.046	N.D.	U	0.21	0.052	0.020	N.D.	U	0.19	0.047	0.024	N.D.	U	0.21	0.054	0.051	N.D.	U	0.19	0.048	0.024
	Fluoranthene	1,500	130	N.D.	Ü	0.19	0.049	0.045	N.D.	Ü	0.21	0.052	0.049	N.D.	Ü	0.19	0.047	0.044	N.D.	Ü	0.21	0.054	0.050	N.D.	Ü	0.19	0.048	0.045
	Fluorene	240	950	N.D.	U	0.19	0.049	0.041	N.D.	U	0.21	0.052	0.044	N.D.	U	0.19	0.047	0.040	N.D.	U	0.21	0.054	0.046	N.D.	U	0.19	0.048	0.041
	Indeno[1,2,3-cd]pyrene	0.092	0.095	N.D.	U	0.19	0.049	0.021	N.D.	U	0.21	0.052	0.022	N.D.	U	0.19	0.047	0.020	N.D.	U	0.21	0.054	0.023	N.D.	U	0.19	0.048	0.020
	1-Methylnaphthalene	4.7	10	N.D.	U	0.19	0.097	0.050	N.D.	U	0.21	0.10	0.054	59		3.8	1.9	0.98	N.D.	U	0.21	0.011	0.055	N.D.	U	0.19	0.096	0.050
	2-Methylnaphthalene	24	10	N.D.	U	0.19	0.049	0.045	N.D.	U	0.21	0.052	0.048	43		3.8	0.95	0.88	N.D.	U	0.21	0.054	0.050	N.D.	U	0.19	0.048	0.045
	Naphthalene Phenanthrene	17 240	21 410	N.D. N.D.	U	0.19 0.19	0.049 0.049	0.033	N.D. N.D.	U	0.21	0.052 0.052	0.035 0.028	140 N.D.	U	3.8 0.19	0.95 0.047	0.64	N.D. N.D.	U	0.21 0.21	0.054 0.054	0.036 0.029	N.D. N.D.	U	0.19 0.19	0.048	0.033 0.026
	Pyrene	180	68	N.D.	II	0.19	0.049	0.026	N.D.	U	0.21	0.052	0.028	N.D.	U	0.19	0.047	0.026 0.019	N.D.	U	0.21	0.054	0.029	N.D.	U	0.19	0.048	0.026
	1,1,1,2-Tetrachloroethane	0.52	50,000	N.D.	U	1	0.049	0.020	N.D.	U	1	0.052	0.021	N.D.	Ü	1	0.047	0.019	N.D.	U	1	0.054	0.022	N.D.	U	1	0.046	0.020
	1,1,2,2-Tetrachloroethane ¹	0.067	500	N.D.	Ü	1	0.5	0.41	N.D.	U	1	0.5	0.41	N.D.	U	1	0.5	0.41	N.D.	U	1	0.5	0.41	N.D.	U	1	0.5	0.41
	1,1,1-Trichloroethane	200	970	N.D.	U	5	0.5	0.3	N.D.	U	5	0.5	0.3	N.D.	U	5	0.5	0.3	N.D.	U	5	0.5	0.3	N.D.	U	5	0.5	0.3
	1,1,2-Trichloroethane	5	50,000	N.D.	U	1	0.5	0.38	N.D.	U	1	0.5	0.38	N.D.	U	1	0.5	0.38	N.D.	U	1	0.5	0.38	N.D.	U	1	0.5	0.38
	1,1-Dichloroethane	2.4	50,000	N.D.	U	5	0.5	0.28	N.D.	U	5	0.5	0.28	N.D.	U	5	0.5	0.28	N.D.	U	5	0.5	0.28	N.D.	U	5	0.5	0.28
	1,1-Dichloroethylene	7	1,500	N.D.	U	1 -	0.5	0.43	N.D.	U	1	0.5	0.43	N.D.	U	1	0.5	0.43	N.D.	U	1	0.5	0.43	N.D.	U	1	0.5	0.43
	1,2,3-Trichloropropane 1,2,4-Trichlorobenzene	0.6 70	50,000 3,000	N.D. N.D.	U	5	1	0.64 0.5	N.D. N.D.	U	5 5	1	0.64	N.D.	U	5 5	1 1	0.64	N.D.	U	5	1	0.64	N.D.	U	5 5	1	0.64
	1,2-Dibromo-3- chloropropane ¹	0.04	10	N.D.	U	10	2	1.2	N.D.	U	10	2	1.2	N.D.	U	10	2	1.2	N.D.	U	10	2	1.2	N.D.	U	10	2	1.2
	1,2-Dibromoethane ¹	0.04	50,000	N.D.	Ü	1	0.5	0.24	N.D.	U	1	0.5	0.24	N.D.	Ü	1	0.5	0.24	N.D.	U	1	0.5	0.24	N.D.	Ü	1	0.5	0.24
	1,2-Dichlorobenzene	600	10	N.D.	Ü	1	0.5	0.46	N.D.	Ü	1	0.5	0.46	N.D.	Ü	1	0.5	0.46	N.D.	Ü	1	0.5	0.46	N.D.	Ü	1	0.5	0.46
	1,2-Dichloroethane ¹	0.15	7,000	N.D.	U	1	0.5	0.24	N.D.	U,IH	1	0.5	0.24	N.D.	U,IH	1	0.5	0.24	N.D.	U,IH	1	0.5	0.24	N.D.	U	1	0.5	0.24
	1,2-Dichloropropane	5	10	N.D.	U	5	0.5	0.42	N.D.	U	5	0.5	0.42	N.D.	U	5	0.5	0.42	N.D.	U	5	0.5	0.42	N.D.	U	5	0.5	0.42
	1,3-Dichlorobenzene	180	5	N.D.	U	1	0.5	0.4	N.D.	U	1	0.5	0.4	N.D.	U	1	0.5	0.4	N.D.	U	1	0.5	0.4	N.D.	U	1	0.5	0.4
	1,3-Dichloropropene (total of cis/trans)	0.43	50,000	N.D.	U	1	0.5	0.25	N.D.	U	1	0.5	0.25	N.D.	U	1	0.5	0.25	N.D.	U	1	0.5	0.25	N.D.	U	1	0.5	0.25
	1,4-Dichlorobenzene	75	5	N.D.	U	1	0.5	0.43	N.D.	U	1	0.5	0.43	N.D.	U	1	0.5	0.43	N.D.	U	1	0.5	0.43	N.D.	U	1	0.5	0.43
	Acetone	22,000	20,000	N.D.	U,ICH	20	10	6	N.D.	U,ICH	20	10	6	N.D.	U,ICH	20	10	6	N.D.	U,ICH	20	10	6	N.D.	U,ICH	20	10	6
	Benzene Bromodichloromethane ¹	5 0.12	170 50,000	N.D. N.D.	U	5	0.5 0.5	0.14 0.21	N.D. N.D.	U	5	0.5 0.5	0.14	N.D. N.D.	U	5	0.5 0.5	0.14 0.21	N.D.	U	1 5	0.5 0.5	0.14 0.21	N.D. N.D.	U	5	0.5	0.14 0.21
	Bromoform	80	510	N.D.	U	10	1	0.5	N.D.	U	10	1	0.5	N.D.	U	10	1	0.5	N.D.	U	10	1	0.5	N.D.	U	10	1	0.5
EPA 8260B	Bromomethane	8.7	50,000	N.D.	Ü	20	5	3.9	N.D.	U	20	5	3.9	N.D.	Ü	20	5	3.9	N.D.	U	20	5	3.9	N.D.	Ü	20	5	3.9
EPA 6200B	Carbon Tetrachloride	5	520	N.D.	U	1	0.5	0.23	N.D.	U	1	0.5	0.23	N.D.	U	1	0.5	0.23	N.D.	U	1	0.5	0.23	N.D.	U	1	0.5	0.23
	Chlorobenzene	100	50	N.D.	U	5	0.5	0.17	N.D.	U	5	0.5	0.17	N.D.	U	5	0.5	0.17	N.D.	U	5	0.5	0.17	N.D.	U	5	0.5	0.17
	Chloroethane	21,000	16	N.D.	U	10	5	2.3	N.D.	U	10	5	2.3	N.D.	U	10	5	2.3	N.D.	U	10	5	2.3	N.D.	U	10	5	2.3
	Chloroform	70	2,400	N.D.	U	5	0.5	0.46	N.D.	U	5	0.5	0.46	N.D.	U	5	0.5	0.46	N.D.	U	5	0.5	0.46	N.D.	U	5	0.5	0.46
	Chloromethane cis 1.2 Dichloroethylene	1.8 70	50,000 50,000	N.D.	U, IJ U	10	2	1.8	N.D.	U, IJ	10	0.5	1.8	N.D. N.D.	U, IJ	10	2	1.8	N.D.	U, IJ	10	2	1.8	N.D. N.D.	U, IJ U	10 1	2	1.8 0.48
	cis-1,2-Dichloroethylene Dibromochloromethane ¹	0.16	50,000	N.D.	U	1	0.5 0.5	0.48 0.25	N.D. N.D.	U	1	0.5	0.48	N.D.	U	1	0.5	0.48 0.25	N.D.	U	1	0.5 0.5	0.48 0.25	N.D.	U	1	0.5	0.48
	Ethylbenzene	700	30	N.D.	U	1	0.5	0.23	N.D.	U	1	0.5	0.25	0.15	J	1	0.5	0.23	N.D.	U	1	0.5	0.23	N.D.	U	1	0.5	0.23
	Hexachlorobutadiene	0.86	6	N.D.	Ű	1	0.5	0.32	N.D.	U	1	0.5	0.32	N.D.	Ŭ	1	0.5	0.32	N.D.	U	1	0.5	0.32	N.D.	Ü	1	0.5	0.32
	Methyl ethyl ketone (2- Butanone)	7,100	8,400	N.D.	U	10	5.0	2.2	N.D.	U,ICH	10	5.0	2.2	N.D.	U,ICH	10	5.0	2.2	N.D.	U,ICH	10	5.0	2.2	N.D.	U	10	5.0	2.2
	Methyl isobutyl ketone (4- Methyl-2-Pentanone)	2,000	1300	N.D.	U	10	5.0	4.4	N.D.	U	10	5.0	4.4	N.D.	U	10	5.0	4.4	N.D.	U	10	5.0	4.4	N.D.	U	10	5.0	4.4
	Methyl tert-butyl Ether	12	5	N.D.	U	1	0.5	0.31	N.D.	U	1	0.5	0.31	N.D.	U	1	0.5	0.31	N.D.	U	1	0.5	0.31	N.D.	U	1	0.5	0.31
	Methylene chloride	4.8	9,100	N.D.	U	5	1.0	0.64	N.D.	U	5	1.0	0.64	N.D.	U	5	1.0	0.64	N.D.	U	5	1.0	0.64	N.D.	U	5	1.0	0.64
	Styrene	100	10	N.D.	U	1	0.5	0.17	N.D.	U	1 5	0.5	0.17	N.D.	U	1 5	0.5	0.17	N.D.	U	1 -	0.5	0.17		U	1	0.5	0.17
	Tetrachloroethylene	5 1,000	170 40	N.D.	U	5	0.5	0.39	N.D.	U	5 1	0.5	0.39	N.D. N.D.	U	5 1	0.5	0.39	N.D.	U	5 1	0.5	0.39	N.D.	U	5 1	0.5	0.39
	Toluene trans-1.2- Dichloroethylene	1,000	40 260	N.D. N.D.	U	1	0.5 0.5	0.24	N.D. N.D.	U	1	0.5 0.5	0.24	N.D.	U	1	0.5 0.5	0.24 0.37	N.D. N.D.	U	1	0.5 0.5	0.24 0.37	N.D. N.D.	U	1	0.5	0.24
	Trichloroethylene	5	310	N.D.	U	1	0.5	0.37	N.D.	U	1	0.5	0.37	N.D.	U	1	0.5	0.37	N.D.	U	1	0.5	0.37	N.D.	U	1	0.5	0.37
	Vinyl chloride	2	3,400	N.D.	Ü	1	0.5	0.3	N.D.	U	1	0.5	0.3	N.D.	U	1	0.5	0.3	N.D.	U	1	0.5	0.3	N.D.	U	1	0.5	0.3
	Xylenes	10,000	20	N.D.	U	11	1.5	0.23	N.D.	U	1	0.5	0.23	0.32	J	1	0.5	0.23	N.D.	U	1	0.5	0.23	N.D.	U	11	1.5	0.23
EPA 6020	Dissolved Lead	15	50,000	-	-	-	-	-	0.0976	J	1	0.2	0.0898	N.D.	U	1	0.2	0.0898	N.D.	U	1	0.2	0.0898	N.D.	U	1	0.2	0.0898
EPA 200.8	Total Lead	15	50,000	0.211	J	1.0	-	0.0898	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-

The data are in micrograms per liter (μ g/L). Shaded values exceeded the DOH EALs.

1 LOD for this analyte exceeds the DOH EAL

- Not Analyzed

DL DOH EALS EPA HD ICH ICJ

Not Analyzed
This sample was analyzed by EPA Method 200.8 and therefore does not have an LOD
Detection Limit or Method Detection Limit (MDL)
DOH Tier 1 Environmental Action Levels for groundwater where groundwater is a current drinking water source and surface water is greater than 150 meters from the site (DOH, Fall 2011).
Environmental Protection Agency
The chromatographic pattern was inconsistent with the profile of the reference fuel standard.
Initial calibration verification recovery is above the control limit for this analyte.
Initial calibration verification recovery is below the control limit for this analyte.

Analyte was detected at a concentration below the LOQ and above the DL. Reported value is estimated. Limit of Detection Limit of Quantitation
Not Detected Qualifiers
Total Petroleum Hydrocarbons as diesel

J LOD LOQ N.D. Q TPH-d TPH-g U

Total Petroleum Hydrocarbons as gasoline
Undetected at DL and is reported as less than the LOD.

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SECTION 3 – DATA QUALITY ASSESSMENT

A data quality assessment, which consists of a review of the overall groundwater sample collection and analysis process, was performed in order to determine whether the analytical data generated met the quality objectives for the project. The data quality assessment was performed in accordance with the approved WP/SAP (ESI, 2012). The field quality control program consisted of standardized sample collection and management procedures, and the collection of field duplicate samples, matrix spike samples, and trip blank samples. The laboratory quality assurance program consisted of the use of standard analytical methods and the preparation and analyses of Matrix Spike [MS]/Matrix Spike Duplicate [MSD] samples, surrogate spikes, blanks, and Laboratory Control Samples [LCSs]/Laboratory Control Sample Duplicates [LCSDs].

3.1 DATA VALIDATION AND ASSESSMENT

The objective of data validation is to provide data of known quality for project decisions. Data quality is judged in terms of Precision, Accuracy, Representativeness, Completeness, Comparability, and Sensitivity [PARCCS]. A number of factors may affect the quality of data, including: sample collection methods, sample analysis methods, and adherence to established procedures for sample collection, preservation, management, shipment, and analysis.

Precision

Precision is defined as the reproducibility of replicate measurements. Precision is evaluated by Relative Percentage Difference [RPD] of field duplicates, LCS/LCSD, and MS/MSD results. Field duplicate and MS/MSD samples were collected at a rate of approximately 10% of primary samples. Field duplicates were sent to the laboratory along with the primary samples.

The RPDs of detected analytes for the primary and field duplicate samples (ES114 and ES115) are provided in Table 3.1. A precision of less than 50% for duplicate pairs is required by the DON Project Procedures Manual to be considered acceptable (DON 2007). All duplicate RPDs are less than the acceptable maximum, except for lead. An RPD of 200% was assigned to the ES114 and ES115 (RHMW02) lead results because lead was detected in only one of the replicate samples from the duplicate pair. The lead result for sample ES115 was below the limit of quantitation [LOQ]; consequently, the RPD exceedance signified the anticipated decrease in precision below the LOQ, but was not indicative of a QC issue. Therefore, it is unlikely to have an effect on data usability. In addition, all RPDs for MS/MSD and LCS/LCSD pairs were also within the control limits, and the data precision is considered acceptable.

Accuracy

Accuracy is defined as the degree of conformity of a measurement to a standard or true value. Accuracy is evaluated through measurement of the percent recovery of an analyte in a reference standard or spiked sample. Accuracy limits for surrogates, laboratory control spike, MS, and MSD samples are either prescribed by the Department of Defense [DoD] or

established by the individual laboratory. The acceptance criteria for accuracy are dependent on the analytical method and are based on historical laboratory or DoD data.

Between July 2006 and July 2010, naphthalene was analyzed for by both EPA Methods 8260B and 8270C, and both results were reported. In September 2005 and in all data beginning in October 2010, only results using EPA Method 8270C were reported. Naphthalene has historically only been detected at concentrations above the DOH EALs in well RHMW02. In this well, concentrations of naphthalene detected in each sample by EPA Method 8260B were generally two to three times higher than those detected by EPA Method 8270C. We assume this is due to the better preservation of VOCs associated with the use of EPA Method 8260B. This suggests that the naphthalene results provided by EPA Method 8270C may be biased low. Since March 2014, naphthalene concentrations in RHMW02 have exceeded DOH EALs for both gross contamination and drinking water toxicity. Therefore, a low bias is unlikely to affect project decisions.

Results for TPH-d in samples ES113 through ES118 were flagged "HD." The laboratory indicated a mismatch between the calibration standard and the TPH-d chromatographic profile. Mismatches of this type are not uncommon. Even though chromatograms are not part of the standard laboratory package, ESI was able to review the chromatograms from RHMW02 dating back to October 2012. The chromatograms of groundwater samples from RHMW02 did not significantly differ between each event, but did not match a standard chromatogram of JP-8 in groundwater.

The MS and MSD recoveries were above the control limits for TPH-d, and the associated sample results may be biased high. The TPH-d results from this sample delivery group (ES113 to ES116) have been flagged "J", estimated, to indicate the increased error caused by possible matrix interference. Naphthalene, 2-methylnaphthalene and 1-methylnaphthalene concentrations for ES114, the primary sample on which the MS/MSD were performed, were significantly higher than the added spike concentration, which prevented an accurate evaluation of the MS/MSD recovery for these analytes.

All other MS/MSD recoveries and all LCS and surrogate spike recoveries were within acceptable recovery limits; therefore, the data accuracy for this monitoring event is considered acceptable.

Representativeness

Representativeness is the degree that data accurately and precisely represents a characteristic of a population, parameter variations at a sampling point, or an environmental condition. Representativeness was achieved by conducting sampling in accordance with the sample collection procedures described in the project WP/SAP, including standardized sample collection methods (ESI, 2012).

Representativeness is also evaluated through the compliance with the standardized sample holding time and sample preservation methods, and through the analysis of blank samples, including method blank and trip blank samples. For this sampling event, all sample holding time and sample preservation were consistent with EPA guidance.

For this sampling event, one trip blank was included in every cooler containing samples for VOC and TPH-g analysis to assess the potential for contamination during sample transport. Two trip blanks were collected. 1,4-Dichlorobenzene was the only COPC detected, and in only one of the trip blanks at a concentration below the LOD; however, 1,4-dichlorobenzene was not detected in any of the groundwater samples, excluding the possibility that project samples were contaminated with 1,4-dichlorobenzene during handling or transport. Based on the assessment of representativeness, the groundwater sample data are considered representative of the groundwater quality on site. The trip blank results are provided in Table 3.1.

Completeness

Completeness is defined as the overall percentage of valid analytical results (including estimated results) compared to the total number of analytical results reported by the analytical laboratory. No data were rejected for this project, and therefore the completeness goal for this project (90%), was successfully met.

Comparability

Comparability expresses the confidence with which one data set can be compared to another data set. Comparability can be related to accuracy and precision because these quantities are measures of data reliability. Data with acceptable precision and accuracy are considered comparable if collection techniques, analytical procedures, methods and reporting are equivalent.

As noted above, between July 2006 and July 2010, naphthalene was analyzed for using both EPA Methods 8260B and 8270C, and in September 2005 and between October 2010 and the most recent event, only results using EPA Method 8270C were reported. In general, EPA Method 8260B resulted in higher, and as discussed above, likely more accurate, results than EPA Method 8270C. However, for the sake of comparability with results from recent events, EPA Method 8270C was used for naphthalene analysis in this event. Consequently, the low bias associated with Method 8270C should be considered when making project decisions.

All project samples for TPH-g analysis through July 2010 were analyzed by EPA Method 8015; beginning in October 2010, EPA Method 8260B was used. There was no event where both methods were used; consequently, there is no way to directly compare the results obtained by method and to assess potential bias. However, there is no reason to believe that using either method should bias the data, and the TPH-g data for all events should be comparable.

Other than the naphthalene bias discussed above, no issues with comparability were identified. The results are considered comparable within this data set and with the data collected from recent sampling events.

Sensitivity

The LOQs are established by the laboratory based on the limits of detection [LODs] or instrument detection limits, historical data, and EPA limits established for the various methods. The LOQs for samples may require adjustment by the laboratory due to matrix interference or if high levels of target analytes necessitate dilution before analysis. Matrix interference and sample dilutions have the effect of decreasing sensitivity and increasing the LOQs. Laboratory LODs and LOQs for several analytes (EPA Methods 8260 and 8270) for this event differed from the LODs and LOQs in the WP/SAP because the laboratory updates them quarterly and in some cases, dilution was necessary due to the presence of high concentrations of analytes.

For this event, LODs and LOQs for several analytes were greater than the DOH EALs (as stated in the WP/SAP), and therefore it would not be possible to detect the analytes at concentrations greater than the DOH EALs but below the LODs. The lack of the required sensitivity should be considered when making project decisions. The affected analytes for this monitoring event are 1,2,3-trichloropropane, 1,2-dibromo-3-chloropropane, 1,2-dibromoethane, 1,2-dichloropropene, bromodichloromethane, chloromethane, dibromochloromethane, 1,1,2,2-tetrachloroethane, and dibenzo[a,h]anthracene.

3.2 DATA ASSESSMENT AND USABILITY CONCLUSIONS

The PARCCS criteria were evaluated, and with a few exceptions, all criteria were met. Based on the high MS/MSD recoveries for TPH-d, it is possible that sample results for TPH-d may be biased high. This increased error should be kept in mind when comparing TPH-d results from this sampling event to those of previous events, but should not significantly affect data usability. Other than this issue, the data assessment concludes that all data generated during this event are usable for the intended purpose.

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TABLE 3.1 Quality Control Results for Groundwater Sampling (October 27 and 28, 2014) Red Hill Bulk Fuel Storage Facility October 2014 Quarterly Monitoring Report

	Chemical Constituent	DOH I		RHM	W02 (ES			Quarter		02 (ES115		-			ES Tr	ip (10/27/	2014)		ES Trip (10/28/2014)					
Method				1		T			1				RPD Duplicate		1		1							
Metriod		Drinking Water Toxicity	Gross Contamination	Result	Q	LOQ	LOD	DL	Result	Q	LOQ	LOD	DL	(%)	Result	Q	LOQ	LOD	DL	Result	Q	LOQ	LOD	DL
EPA 8015B	TPH-d	190	100	2,000	HD,J	25	12	11	2,000	HD,J	25	12	11	0.00	-	-	-	-	-	-	-	-	-	-
EPA 8260B	TPH-g	100 370	100 20	57 0.53		50 0.19	30 0.047	26 0.025	53 0.53	J	50 0.19	30 0.047	26 0.025	7.27 0.00	N.D.	U -	50	30	26	N.D.	U	50	30	26
	Acenaphthylene Acenaphthylene	240	2,000	0.53 N.D.	U	0.19	0.047	0.025	0.53 N.D.	U	0.19	0.047	0.025	NA	-	-	-	-	-	-	-	-	-	-
	Anthracene	1,800	22	N.D.	Ü	0.19	0.047	0.042	N.D.	U	0.19	0.047	0.042	NA NA	_	-	-	-	-	-	-	-	-	-
	Benzo[a]anthracene	0.092	4.7	N.D.	U	0.19	0.047	0.031	N.D.	U	0.19	0.047	0.031	NA	-	-	-	-	-	-	-	-	-	-
	Benzo[g,h,i]perylene	1,500	0.13	N.D.	U	0.19	0.095	0.078	N.D.	U	0.19	0.095	0.078	NA	-	-	-	-	-	-	-	-	-	-
	Benzo[a]pyrene	0.2	0.81	N.D.	U	0.19	0.047	0.021	N.D.	U	0.19	0.047	0.021	NA	-	-	-	-	-	-	-	-	-	-
	Benzo[b]fluoranthene	0.092	0.75	N.D.	U	0.19	0.047	0.017	N.D.	U	0.19	0.047	0.017	NA NA	-	-	-	-	-	-	-	-	-	-
	Benzo[k]fluoranthene	0.92 9.2	0.4	N.D.	U	0.19 0.19	0.047 0.047	0.029 0.024	N.D. N.D.	U	0.19 0.19	0.047 0.047	0.029 0.023	NA NA	-	-	-	+	-	-	-	-	-	-
EPA 8270C	Chrysene Dibenzoſa.hlanthracene	0.0092	0.52	N.D.	U	0.19	0.047	0.024	N.D.	U	0.19	0.047	0.023	NA NA	-	-	-	-	-	-	-	-	-	-
	Fluoranthene	1,500	130	N.D.	Ü	0.19	0.047	0.044	N.D.	U	0.19	0.047	0.043	NA NA	_	-	_	_	_	-	_		_	_
	Fluorene	240	950	N.D.	Ü	0.19	0.047	0.040	N.D.	Ü	0.19	0.047	0.040	NA	-	-	-	-	-	-	-	-	-	-
	Indeno[1,2,3-cd]pyrene	0.092	0.095	N.D.	U	0.19	0.047	0.020	N.D.	U	0.19	0.047	0.020	NA	-	-	-	-	-	-	-		-	-
	1-Methylnaphthalene	4.7	10	59		3.8	1.9	0.98	54		3.8	1.9	0.98	8.85	-	-	-	-	-	-	-	-	-	-
	2-Methylnaphthalene	24	10	43		3.8	0.95	0.88	36		3.8	0.95	0.88	17.72	-	-	-	-	-	-	-	-	-	-
	Naphthalene	17	21	140	L	3.8	0.95	0.64	130		3.8	0.95	0.64	7.41	-	-	-	-	-	-	-	-	-	-
	Phenanthrene Pyrene	240 180	410 68	N.D. N.D.	U	0.19 0.19	0.047 0.047	0.026 0.019	N.D. N.D.	U	0.19 0.19	0.047 0.047	0.026 0.019	NA NA	-	-	-	-	-	-	-	-	-	-
	1,1,1,2-Tetrachloroethane	0.52	50,000	N.D.	U	1	0.047	0.019	N.D.	U	1	0.047	0.019	NA NA	N.D.	U	1	0.5	0.4	N.D.	U	1	0.5	0.4
	1,1,2,2-Tetrachloroethane	0.067	500	N.D.	Ü	1	0.5	0.41	N.D.	U	1	0.5	0.41	NA	N.D.	Ü	1	0.5	0.41	N.D.	Ü	1	0.5	0.41
	1,1,1-Trichloroethane	200	970	N.D.	U	5	0.5	0.3	N.D.	U	5	0.5	0.3	NA	N.D.	U	5	0.5	0.3	N.D.	U	5	0.5	0.3
	1,1,2-Trichloroethane	5	50,000	N.D.	U	1	0.5	0.38	N.D.	U	1	0.5	0.38	NA	N.D.	U	1	0.5	0.38	N.D.	U	1	0.5	0.38
	1,1-Dichloroethane	2.4	50,000	N.D.	U	5	0.5	0.28	N.D.	U	5	0.5	0.28	NA	N.D.	U	5	0.5	0.28	N.D.	U	5	0.5	0.28
	1,1-Dichloroethylene	7	1,500	N.D.	U	1	0.5	0.43	N.D.	U	1	0.5	0.43	NA NA	N.D.	U	1 -	0.5	0.43	N.D.	U	1	0.5	0.43
	1,2,3-Trichloropropane 1,2,4-Trichlorobenzene	0.6 70	50,000 3,000	N.D. N.D.	U	5 5	1	0.64 0.5	N.D. N.D.	U	5	1	0.64	NA NA	N.D. N.D.	U	5 5	1	0.64	N.D. N.D.	U	5	1	0.64 0.5
	1,2-Dibromo-3- chloropropane	0.04	10	N.D.	U	10	2	1.2	N.D.	U	10	2	1.2	NA NA	N.D.	U	10	2	1.2	N.D.	U	10	2	1.2
	1,2-Dibromoethane	0.04	50,000	N.D.	Ü	1	0.5	0.24	N.D.	U	1	0.5	0.24	NA	N.D.	Ü	1	0.5	0.24	N.D.	Ü	1	0.5	0.24
	1,2-Dichlorobenzene	600	10	N.D.	U	1	0.5	0.46	N.D.	U	1	0.5	0.46	NA	N.D.	U	1	0.5	0.46	N.D.	U	1	0.5	0.46
	1,2-Dichloroethane	0.15	7,000	N.D.	U,IH	1	0.5	0.24	N.D.	U,IH	1	0.5	0.24	NA	N.D.	U,IH	1	0.5	0.24	N.D.	U	1	0.5	0.24
	1,2-Dichloropropane	5	10	N.D.	U	5	0.5	0.42	N.D.	U	5	0.5	0.42	NA	N.D.	U	5	0.5	0.42	N.D.	U	5	0.5	0.42
	1,3-Dichlorobenzene	180	5	N.D.	U	1	0.5	0.4	N.D.	U	1	0.5	0.4	NA NA	N.D.	U	1	0.5	0.4	N.D.	U	1	0.5	0.4
	1,3-Dichloropropene (total of cis/trans) 1,4-Dichlorobenzene	0.43 75	50,000 5	N.D. N.D.	U	1	0.5 0.5	0.25 0.43	N.D. N.D.	U	1	0.5 0.5	0.25 0.43	NA NA	N.D. N.D.	U	1	0.5 0.5	0.25 0.43	N.D. 0.43	U	1	0.5 0.5	0.25 0.43
	Acetone	22,000	20,000	N.D.	U.ICH	20	10	6	N.D.	U,IC	20	10	6	NA NA	N.D.	U,IC	20	10	6	0.43 N.D.	U,IC	20	10	6
	Benzene	5	170	N.D.	U	1	0.5	0.14	N.D.	U	1	0.5	0.14	NA NA	N.D.	U	1	0.5	0.14	N.D.	U	1	0.5	0.14
	Bromodichloromethane	0.12	50,000	N.D.	Ü	5	0.5	0.21	N.D.	U	5	0.5	0.21	NA	N.D.	Ü	5	0.5	0.21	N.D.	Ü	5	0.5	0.21
	Bromoform	80	510	N.D.	U	10	1	0.5	N.D.	U	10	1	0.5	NA	N.D.	U	10	1	0.5	N.D.	U	10	1	0.5
EPA 8260B	Bromomethane	8.7	50,000	N.D.	U	20	5	3.9	N.D.	U	20	5	3.9	NA	N.D.	U	20	5	3.9	N.D.	U,ICJ	20	5	3.9
	Carbon Tetrachloride	5	520	N.D.	U	1	0.5	0.23	N.D.	U	1	0.5	0.23	NA	N.D.	U	1	0.5	0.23	N.D.	U	1	0.5	0.23
	Chlorothone	100	50	N.D.	U	5	0.5	0.17	N.D.	U	5	0.5 5	0.17	NA NA	N.D.	U	5	0.5	0.17	N.D.	U	5	0.5	0.17
	Chloroethane Chloroform	21,000 70	16 2,400	N.D. N.D.	U	10 5	5 0.5	2.3 0.46	N.D. N.D.	U	10 5	0.5	2.3 0.46	NA NA	N.D. N.D.	U	10 5	5 0.5	2.3 0.46	N.D. N.D.	U	10 5	5 0.5	2.3 0.46
	Chloromethane	1.8	50,000	N.D.	U, IJ	10	2	1.8	N.D.	U, IJ	10	2	1.8	NA NA	N.D.	U, IJ	10	2	1.8	N.D.	U	10	2	1.8
	cis-1,2-Dichloroethylene	70	50,000	N.D.	U	1	0.5	0.48	N.D.	U	1	0.5	0.48	NA	N.D.	U	1	0.5	0.48	N.D.	U	1	0.5	0.48
	Dibromochloromethane	0.16	50,000	N.D.	U	1	0.5	0.25	N.D.	U	1	0.5	0.25	NA	N.D.	U	1	0.5	0.25	N.D.	U	1	0.5	0.25
ĺ	Ethylbenzene	700	30	0.15	J	1	0.5	0.14	0.14	J	1	0.5	0.14	6.90	N.D.	U	1	0.5	0.14	N.D.	U	1	0.5	0.14
ĺ	Hexachlorobutadiene	0.86	6	N.D.	U	1	0.5	0.32	N.D.	U	1	0.5	0.32	NA NA	N.D.	U	1	0.5	0.32	N.D.	U	1	0.5	0.32
ĺ	Methyl ethyl ketone (2-Butanone)	7,100 2,000	8,400 1300	N.D.	U,ICH	10	5.0	2.2	N.D.	U,IC	10	5.0	2.2	NA NA	N.D.	U,IC	10	5.0	2.2	N.D.	U	10	5.0	2.2
ĺ	Methyl isobutyl ketone (4-Methyl-2-Pentanone) Methyl tert-butyl Ether	2,000	1300	N.D. N.D.	U	10	5.0 0.5	4.4 0.31	N.D. N.D.	U	10 1	5.0 0.5	4.4 0.31	NA NA	N.D. N.D.	U	10	5.0 0.5	4.4 0.31	N.D. N.D.	U	10 1	5.0 0.5	4.4 0.31
Ī	Methylene chloride	4.8	9,100	N.D.	U	5	1.0	0.64	N.D.	U		1.0	0.64	NA NA	N.D.	U	5	1.0	0.64	N.D.	U	5	1.0	0.64
Ī	Styrene	100	10	N.D.	U	1	0.5	0.17	N.D.	U	1	0.5	0.17	NA	N.D.	U	1	0.5	0.17	N.D.	U	1	0.5	0.17
ĺ	Tetrachloroethylene	5	170	N.D.	Ü	5	0.5	0.39	N.D.	U	5	0.5	0.39	NA	N.D.	Ü	5	0.5	0.39	N.D.	Ü	5	0.5	0.39
ĺ	Toluene	1,000	40	N.D.	U	1	0.5	0.24	N.D.	U	1	0.5	0.24	NA	N.D.	U	1	0.5	0.24	N.D.	U	1	0.5	0.24
	trans-1,2- Dichloroethylene	100	260	N.D.	U	1	0.5	0.37	N.D.	U	1	0.5	0.37	NA	N.D.	U	1	0.5	0.37	N.D.	U	1	0.5	0.37
	Trichloroethylene	5	310	N.D.	U	1	0.5	0.37	N.D.	U	1	0.5	0.37	NA NA	N.D.	U	1	0.5	0.37	N.D.	U	1	0.5	0.37
Ī	Vinyl chloride Xylenes	10,000	3,400 20	N.D. 0.32	U J	1	0.5 0.5	0.3 0.23	N.D. 0.29	J	1	0.5 0.5	0.3 0.23	NA 9.84	N.D. N.D.	U	1 11	0.5 1.5	0.3 0.23	N.D. N.D.	U	1 11	0.5 1.5	0.3 0.23
EPA 6020	Lead	10,000	50,000	0.32 N.D.	U		0.5	0.23	0.29			0.5		9.84	N.D.	-	- 11	1.5	0.23	IN.D.	Ū	- 11	1.5	0.23
EFA 0020	I LCau	ded the DOH EALs	30.000	IV.D.	U		0.2	0.0090	U. 105	J		U.Z	0.0090	200		<u> </u>	-	<u> </u>		-		-		

The data are in micrograms per liter (µg/L). Shaded values exceeded the DOH EALs.

Not Analyzed

DOH EALs

DOH Tier 1 Environmental Action Levels for groundwater where groundwater is a current drinking water source and surface water is greater than 150 meters from the site (DOH, Fall 2011).

Detection Limit or Method Detection Limit (MDL)

EPA

Environmental Protection Agency

The detarance fine later darks.

DOH EALS
DL
EPA
HD
ICH
ICJ
J

Environmental Protection Agency
The chromatographic pattern was inconsistent with the profile of the reference fuel standard.
Initial calibration verification recovery is above the control limit for this analyte.
Initial calibration verification recovery is below the control limit for this analyte.
Analyte was detected at a concentration below the LOQ and above the DL. Reported value is estimated.

LOD LOQ NA N.D. Q TPH-g TPH-d U

Limit of Detection Limit of Quantitation Both results for duplicate pair were non-detect, no RPD calculations

Not Detected

Not Detected
Qualifiers
Total Petroleum Hydrocarbons as gasoline
Total Petroleum Hydrocarbons as diesel
Undetected at DL and is reported as less than the LOD.

Contract No. N62742-12-D-1853

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SECTION 4 – SUMMARY, CONCLUSIONS, AND RECOMMENDATIONS

On October 27 and 28, 2014, ESI personnel collected groundwater samples from four monitoring wells at the RHSF (wells RHMW01, RHMW02, RHMW03, and RHMW05) and one sampling point at Red Hill Shaft (RHMW2254-01).

The groundwater sampling was conducted as part of the long-term groundwater and soil vapor monitoring program at the RHSF, under NAVFAC Contract Number N62742-12-D-1853. The sampling was conducted in accordance with the approved WP/SAP prepared by ESI. A summary of the analytical results is provided below.

- RHMW01 TPH-d (120 μg/L) and lead (0.0976 μg/L) were the only analytes detected. The
 TPH-d concentration exceeded the DOH EAL for gross contamination, but was below the
 SSRBL of 4,500 μg/L for TPH. None of the other chemical constituents analyzed for were
 detected at concentrations above their respective DOH EAL.
- RHMW02 TPH-d (2,000 μg/L), TPH-g (57 and 53 μg/L), xylenes (0.32 and 0.29 μg/L), acenaphthene (0.53 μg/L), ethylbenzene (0.15 and 0.14 μg/L), 1-methylnaphthalene (59 and 54 μg/L), 2-methylnaphthalene (43 and 36 μg/L), and naphthalene (140 and 130 μg/L) were detected in both the primary and duplicate samples collected. Lead (0.165 μg/L) was only detected in the duplicate sample. TPH-d, 1-methylnaphthalene, 2-methylnaphthalene, and naphthalene were detected at concentrations above their respective DOH EALs for both drinking water toxicity and gross contamination. However, the TPH-d concentrations did not exceed the SSRBL.
- RHMW03 TPH-d (80 μg/L) was the only analyte detected. The concentration did not exceed the DOH EAL or the SSRBL.
- RHMW05 TPH-d (16 μg/L) was the only analyte detected. The concentration did not exceed the DOH EAL or the SSRBL.
- RHMW2254-01 TPH-d (22 μg/L) and total lead (0.211 μg/L) were the only analytes detected. The concentrations did not exceed their respective DOH EAL.

Groundwater Contaminant Trends

- RHMW01 COPCs detected during this round of quarterly sampling are consistent with the
 historical data for RHMW01. TPH-d has historically been detected at concentrations above
 the DOH EAL for both drinking water toxicity and gross contamination. TPH-d
 concentrations have shown an overall decreasing trend from a high of 1,500 μg/L in
 February 2005. The TPH-d concentration exceeded the DOH EAL for gross contamination
 for the first time since April 7, 2014.
- RHMW02 COPCs detected during this round of quarterly sampling are consistent with the
 historical data for RHMW02. TPH-g, TPH-d, 1-methylnaphthalene, 2-methylnaphthalene,
 and naphthalene have historically been detected at concentrations above the DOH EALs.
 During the October 2014 event, concentrations of TPH-d, 1-methylnaphthalene,

2-methylnaphthalene, and naphthalene approximately doubled from the previous event in July 2014 and remained above DOH EALs. The concentrations of 1-methylnaphthalene, 2-methylnaphthalene, and naphthalene increased to their highest levels in over 5 years. The concentrations of these PAHs are also significantly higher than they were in January 2014, when a high TPH-d spike was observed in this well. The concentrations of TPH-g remained below the DOH EALs for gross contamination and drinking water toxicity and were comparable to the concentrations detected during the previous event.

- RHMW03 COPCs detected during this round of quarterly sampling are consistent with the historical data for RHMW03. TPH-d has historically been detected at concentrations above the DOH EALs; however, it has not been detected at concentrations above the DOH EALs since October 2010.
- RHMW05 COPCs detected during this round of quarterly sampling are consistent with the
 historical data for RHMW05. TPH-d has historically been detected in RHMW05 at
 concentrations above the DOH EAL for both drinking water toxicity and gross contamination;
 however, it has not been detected at concentrations above the DOH EALs since January
 2010.
- RHMW2254-01 COPCs detected during this round of quarterly sampling are consistent
 with the historical data for RHMW2254-01. Although the method reporting limits for TPH-d
 exceeded one or both DOH EALs for drinking water toxicity and gross contamination
 between May 2009 and July 2010, TPH-d was last detected in RHMW2254-01 at a
 concentration above the DOH EAL for gross contamination in January 2008.

Conclusions and Recommendations

During the sampling event conducted on October 27 and 28, 2014, TPH-d in RHMW01 and TPH-d, 1-methylnaphthalene, 2-methylnaphthalene, and naphthalene in RHMW02 were detected at concentrations exceeding the DOH EALs. Groundwater contaminant concentrations in wells RHMW03, RHMW05, and RHMW2254-01 remained at low concentrations and did not change significantly from the previous event, or were not detected.

Concentrations of 1-methylnaphthalene, 2-methylnaphthalene, and naphthalene in RHMW02 have shown an increasing trend since March 5, 2014, but have remained below the historic maximums detected in the well. All other analytical results were generally consistent with historical data.

Based on the groundwater monitoring results and the reported release at Tank 5 in January 2014, continued groundwater monitoring at the wells inside the RHSF tunnel is recommended. The next quarterly event is tentatively scheduled for January 2015.

SECTION 5 – FUTURE WORK

Future work includes the first quarter 2015 groundwater monitoring which is tentatively scheduled for January 2015. A quarterly groundwater monitoring report will be prepared to document the sampling event.

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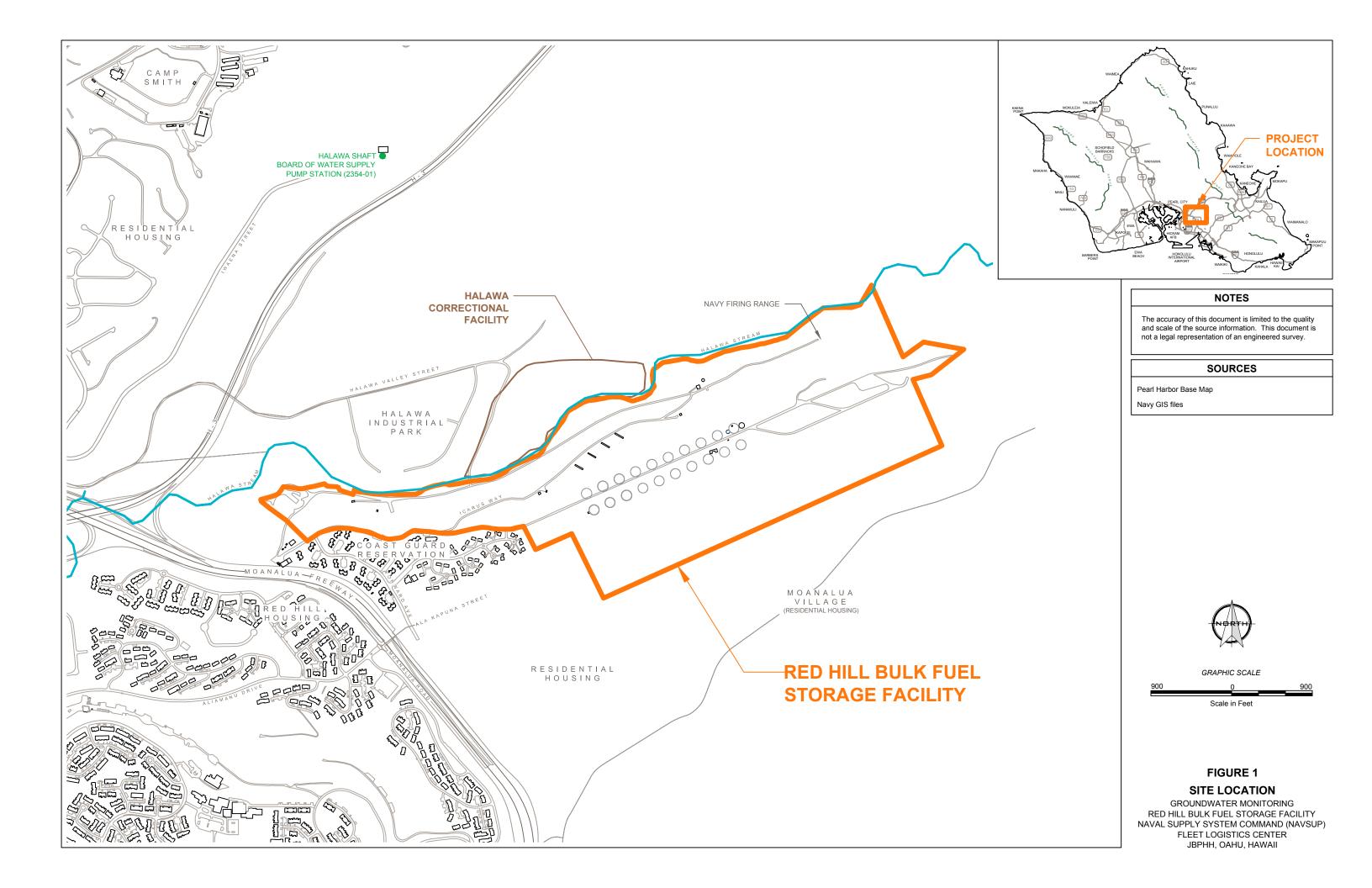
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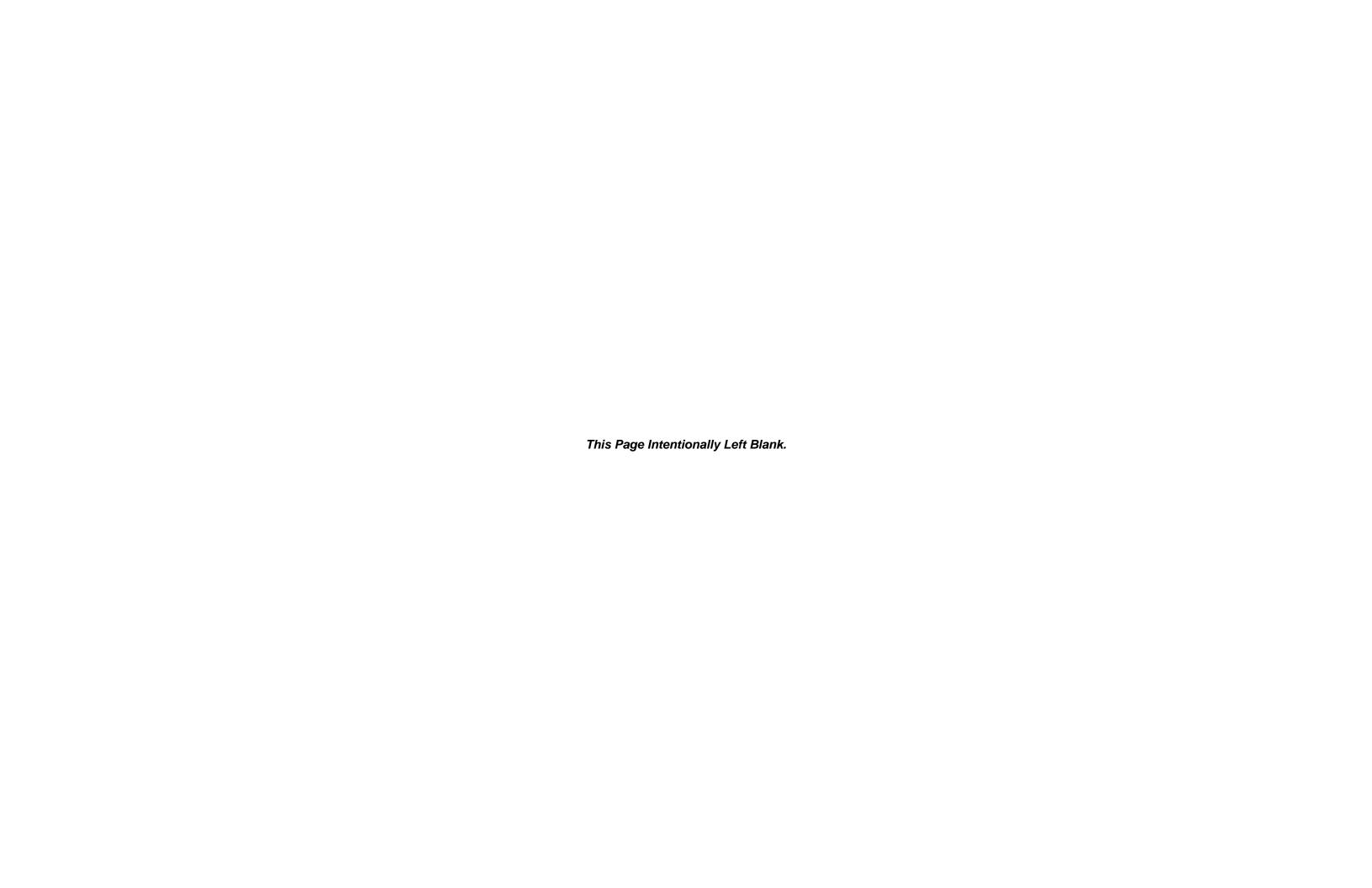
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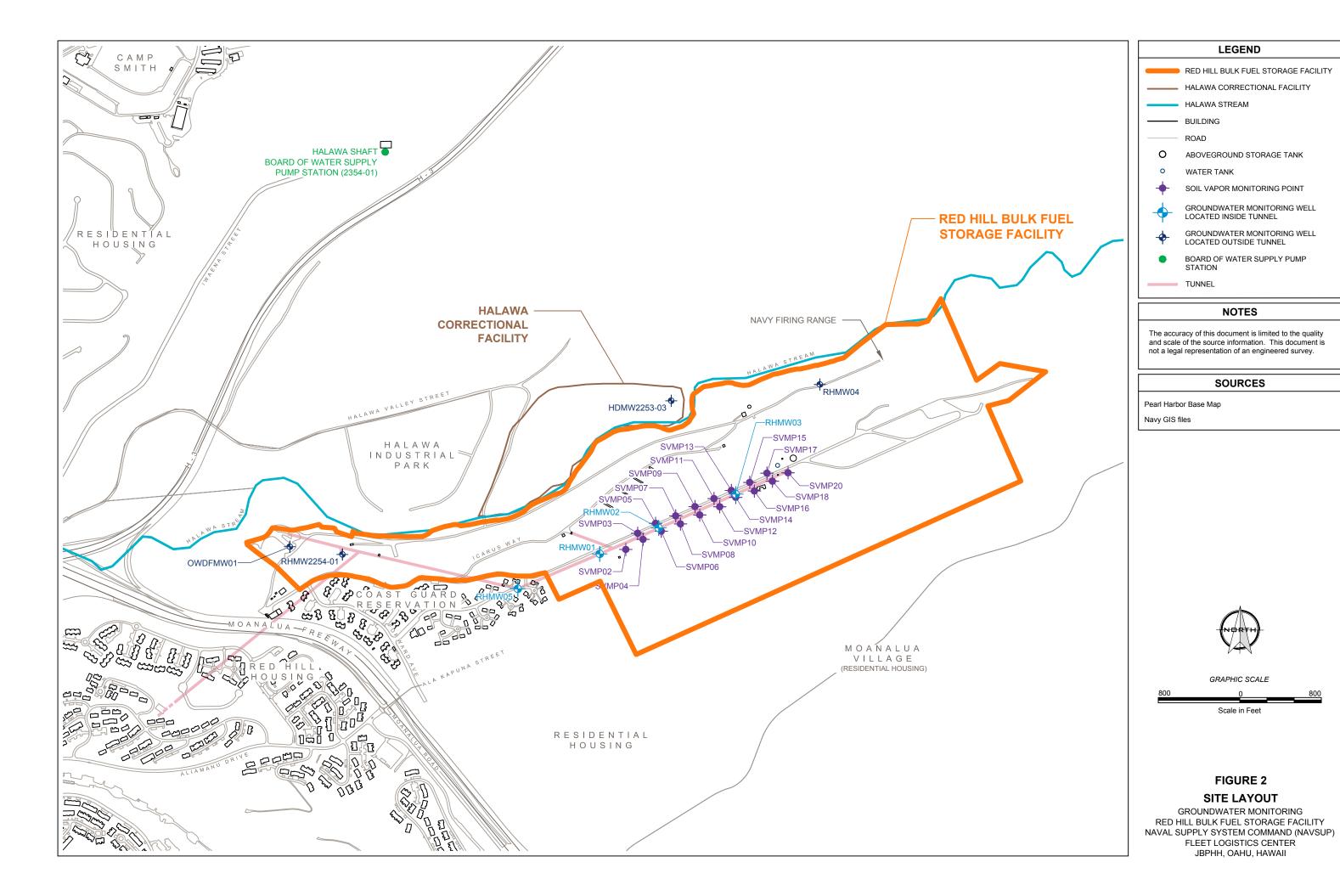
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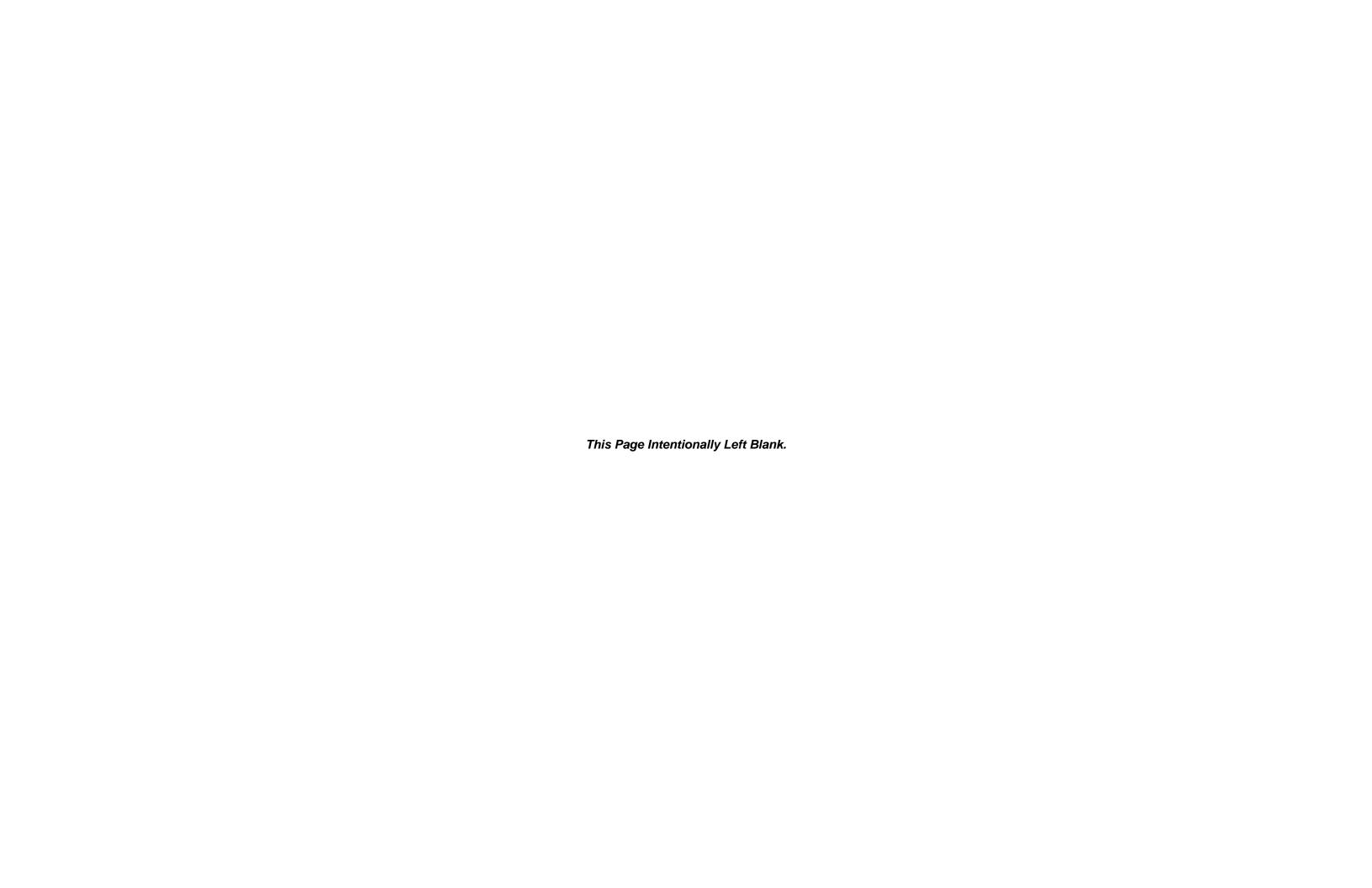
FIGURES











APPENDIX A Groundwater Sampling Logs





Well ID:	RHMW01	Location	on: Red Hill	Bulk Fuel Stora	ige Facility F	Project No.:	112066	
Initial Wa	ater Level:	83.79 ft	Date:	10/27/2014		Гіте: <u>1300</u>)	
Total De	oth of Well:	97.35 ft	Person	nel Involved:	Kirk Markle, Jeff Hattemer			
Length o	ength of Saturated Zone: 13.56 ft		t Weathe	er Conditions:	Not applicable – well is located indoors			
Volume o	of Water to be	e Removed: 3.	5 L Method	of Removal:	Blac	dder Pump		
Water Le	evel After Pur	ging: 83.79	ft Pumpin	ng Rate:	0.19	9 L/min		
Well Pur	ge Data:							
	Volum	е	Conductivity				Redox (ORP)	
Time	Remove	ed pH	(mS/cm)	DO (mg/l)	Temperature	Salinity	(mV)	
1315	0.0 L	8.40	0.342	11.41	25.25		-247.8	
1318	0.5 L	8.39	0.340	2.43	24.88		-289.0	
1322	1.0 L	8.38	0.339	1.64	24.23		-312.0	
1325	1.5 L	8.38	0.335	0.34	23.55	-	-346.0	
1327	2.0 L	8.37	0.335	-0.07	23.41	-	-345.0	
1329	2.5 L	8.35	0.334	-0.06	23.37	-	-342.1	
1331	3.0 L	8.36	0.334	-0.07	23.36	-	-341.2	
1333	3.5 L	8.36	0.332	-0.07	23.34	-	-339.6	
	_							
Sample \	Nithdrawal M	lethod:	Bladder Pur	mp				
Appearar	nce of Sampl	e:						
	Color:		Clear					
	Turbidity	<i>'</i> .	Low	_				
	Sedimer	nt:	None					
	Other:		None					
Laborato	ry Analysis P	arameters and P	reservatives:	TPH-d - 8015;	TPH-g, VOCs - 8	260; PAHs - 82	70c sim;	
				lead - 6020				
Number	and Types of	Sample Containe	ers: <u>6 - 40m</u>	ıl VOAs, 2 - 1L a	ımber jar, 1 - 500n	nl amber jar, 1 -	250ml HDPE	
Sample I	dentification	Numbers: <u>ES</u>	113 [0400]					
Decontar	mination Prod	cedures: Triple	Rinsed					
Notes:	YSI did not h	ave salinity paran	neter.					
Sampled	by: Kirk N	Markle, Jeff Hatte	mer					
•	Delivered to	: Calscienc	e Environment		Transporters: Fed	Ex		
Date: _	10/27/2014				Time: 1530			
			Capacity of Ca	sing (Gallons/Lir	near Feet)			





Well ID: R	HMW02	Location:	Red Hill I	Bulk Fuel Stora	ge Facility P	roject No.:	112066
Initial Water	Level: 86.5	51 ft	Date:	10/27/2014	т	ime: 1011	
Total Depth	of Well:	92.91 ft	Personnel Involved:		Kirk Markle, Jeff Hattemer		
Length of Sa	aturated Zone:	6.40 ft	_ Weather	Conditions:	Not applicable – well is located indoors		
Volume of Water to be Removed: 4.0 L			Method	of Removal:	Blad	der Pump	
Water Level	After Purging:	86.45 ft	_ Pumping	g Rate:	0.17	L/min	
Well Purge							
Time	Volume Removed		onductivity (mS/cm)	DO (mg/l)	Temperature	Salinity	Redox (ORP) (mV)
1029	0.0 L	8.40	0.373	2.40	25.17		291
1035	1.0 L	8.41	0.575	1.30	24.26		298
1040	2.0 L	8.40	0.551	0.97	24.13		-352.4
1045	3.0 L	8.39	0.560	0.82	24.21	-	-349.7
1053	4.0 L	8.35	0.567	0.84	24.25	-	-361.3
•	ndrawal Method	d:	Bladder Pum	р			
Appearance	of Sample: Color:		Clear				
	Turbidity:		Low				
	Sediment:		None				
	Other:		None				
Laboratory A	Analysis Param	eters and Prese	ervatives:	TPH-d - 8015;	TPH-g, VOCs - 82	260; PAHs - 827	0c sim;
				lead - 6020			
Number and	l Types of Sam	ple Containers:	16 - 40ml V	OAs, 6 - 1L am	iber jar, 4 - 500ml	amber jar, 4 - 50	00ml HDPE
Sample Ider	ntification Numb	pers: ES114	[1045], ES11	14 MS/MSD [10	45], ES115 (Dup)	[1115]	
Decontamin	ation Procedure	es: Triple Rins	sed				
		alinity paramete	r.				
Sampled by		0-1 : =		1 -1-	T	=	
Sampled De		Calscience E	nvironmental		Transporters: Fed!	=X	
Date: 10/	27/2014	Car	pacity of Casi	ng (Gallons/Lin	Time: <u>1530</u> ear Feet)		





Well ID: F	RHMW03	Location	: Red Hill	Bulk Fuel Stora	ge Facility	Project No.:	112066
Initial Water	r Level: _	102.78 ft	Date:	10/27/2014		Time: 820	
Total Depth	of Well:	110.12 ft	Personn	Personnel Involved: Kirk Markle, Jeff Hattemer			<u>r</u>
Length of S	aturated 2	Zone: 7.34 ft	Weathe	r Conditions:	Not applicable	- well is located	indoors
Volume of Water to be Removed: 3.0 L			Method	of Removal:	Bla	idder Pump	
Water Level After Purging: 102.88 ft			Pumping	g Rate:	0.	2 L/min	
Well Purge	Data:						
Time	Volum Remov		Conductivity (mS/cm)	DO (mg/l)	Temperature	Salinity	Redox (ORP) (mV)
830	0.0 L		0.823	2.14	26.10	- Saillilly	-206.0
834	1.0 L		0.820	0.95	26.40		-240.1
840	2.0 [0.821	0.87	26.40		-233.1
845	3.0 [0.816	0.71	26.41	-	-238.6
			_	-			
			_				
Sample Wit	hdrawal N	Method:	Bladder Pun	าท			
Appearance			Bladdol 1 dil	<u>h</u>			
. фросполос	Color:		Clear				
	Turbidit	V:	Low				
	Sedime	·	None				
	Other:		None	_			
Labaratan	A b ' F			TDU - 0045	TDU - 1/00-	0000 DALL- 00	701
Laboratory	Anaiysis i	Parameters and Pres	servatives:		TPH-g, VOCs - 8	8260; PAHS - 82	7UC SIM;
Numberon	d Tunaa a	f Cample Containers	o 6 40ml	lead - 6020	mborior 1 F00	Imlambariar 1	250ml UDDE
Sample Ide	• •	of Sample Containers	<u> </u>	VOAS, 2 - 1L a	mber jar, 1 - 500	illi allibei jai, i -	- ZSUIIII HDPE
Decontamir			6 [0900]				
		nave salinity parame					
Sampled by		Markle, Jeff Hatteme					
Sampled Do		,	Environmenta	ıl Lab	Transporters: Fe	dEx	
Date: 10	/27/2014				Time: 1530		
		Ca	apacity of Cas	ing (Gallons/Lir	near Feet)		





Well ID: R	HMW05	Location:	Red Hill	Bulk Fuel Stora	age Facility	Project No.:	112066	
Initial Water	Level: 83.	21 ft	Date:	10/28/2014		Time: 1	013	
Total Depth	of Well: Una	able to Measure	e Personn	Personnel Involved:		Kirk Markle, Jeff Hattemer		
Length of Sa	aturated Zone:	Unknown	Weathe	Weather Conditions:		Not applicable – well is located indoors		
Volume of Water to be Removed: 3.0 L			Method	of Removal:	В	ladder Pump		
Water Level After Purging: 83.21 ft			_ Pumping	g Rate:	0	.23 L/min		
Well Purge								
Time	Volume Removed	pH C	conductivity (mS/cm)	DO (mg/l)	Temperature	e Salinity	Redox (ORP) (mV)	
1405	0.0 L	9.20	0.889	8.83	22.91	<u> </u>	-211.6	
1410	1.0 L	9.61	0.888	8.69	22.87	_	-210.8	
1413	2.0 L	9.22	0.886	8.68	22.87		-210.4	
1418	3.0 L	9.20	0.886	8.67	22.85		-209.1	
						_		
					-			
						_		
					-	_		
Commis Mill	- duarrial Matha		Die delen Dine			_		
•	ndrawal Metho	a:	Bladder Pun	np				
Appearance	Color:		Clear					
	Turbidity:		Clear					
	Sediment:		None					
	Other:		None					
	_							
Laboratory A	Analysis Param	neters and Pres	ervatives:		TPH-g, VOCs -	- 8260; PAHs -	8270c sim;	
				lead - 6020			===	
	• •	nple Containers		I VOAs, 2 - 1L a	amber jar, 1 - 50	00ml amber jar	, 1 - 1L HDPE	
•	ntification Num		8 [1430] ·					
	ation Procedur							
		salinity paramet						
Sampled by Sampled De		e, Jeff Hatteme	er Environmenta	al I ah	Transporters: F	edEv		
•	28/2014	Calsolelice E			Time: 1530	CULX		
10/	Canacity of Casing (Gallons/Linear Feet)							





Well ID:	RHMW2254	<u>-01</u> Locat	on: Red Hil	l Bulk Fuel Stora	age Facility F	Project No.:	112066
Initial Wa	iter Level:	83.44 ft	Date:	10/28/2014	Т	ime: 1110)
Total Dep	oth of Well:	Not applicab	le Person	nel Involved:	Kirk Markl	e, Jeff Hatteme	r
Length of	f Saturated Z	one: Not applic	able Weathe	er Conditions:	Not applicable -	- well is located	indoors
Volume o	of Water to b	e Removed: 5.	0 L Method	d of Removal:	Blad	der Pump	
Water Le	vel After Pur	ging: 83.4	4 ft Pumpir	ng Rate:	0.50) L/min	
Well Purç							
Time	Volum Remov		Conductivity (mS/cm)	DO (mg/l)	Temperature	Salinity	Redox (ORP)
1117	0.0 L	9.60	0.613	DO (mg/l) 13.00	21.97	Samily	(mV) -201.7
1119	1.0 L	9.55	0.606	10.29	21.59		-201.7
1121	2.0 L	9.65	0.603	9.57	21.45		- <u>-200.5</u> -197.5
1123	3.0 L	9.47	0.603	9.43	21.38	 _	-196.2
1125	4.0 L	9.44	0.603	9.43	21.28		-195.0
1127	5.0 L	9.43	0.603	9.17	21.28		-193.7
1127		0.10	0.000	0.17	21.20		100.7
			·				<u> </u>
							<u> </u>
•	Withdrawal M		Bladder Pu	mp			
Appearar	nce of Sampl	e:	0.1				
	Color:		Clear				
	Turbidity		Clear				
	Sedime	nt:	None				
	Other:		None				
Laborato	ry Analysis P	arameters and F	reservatives:	TPH-d - 8015;	; TPH-g, VOCs - 82	260; PAHs - 82 ⁻	70c sim;
				lead - 200.8			
Number a	and Types of	Sample Contain	ers: 6 - 40m	nl VOAs, 2 - 1L a	amber jar, 1 - 500n	nl amber jar, 1 -	· 1L HDPE
Sample I	dentification	Numbers: ES	3117 [1200]			-	
Decontar	mination Prod	cedures: Triple	Rinsed				
Notes: `	YSI did not h	ave salinity para	meter.				
Sampled	by: Kirk N	Markle, Jeff Hatte	emer				
•	Delivered to	: Calscien	ce Environment		Transporters: Fed	Ex	
Date:	10/28/2014		Canacity of Ca	sing (Gallons/Li	Time: 1530		
			Capacity of Ca	ioniy (Galions/Lii	noar reet)		



APPENDIX B Field Notes



Project / Client Date 7 25/14 Personnel 5H 1M

D899 @ Red Hill Sorling meetry

Calibrate PID

A All notes on Field Forms

1500 Depart Red Hill No petroleum olar noted.

in larer france.

Project / Client NAVFAC

Date 10/20/12

73

Purpose SV monsturing GW Sampison

Personnel STIL, KM

0800 Arrive @ 201 Hill, Calibrate

PID, Sterry meeting

DIW: Set up pump.

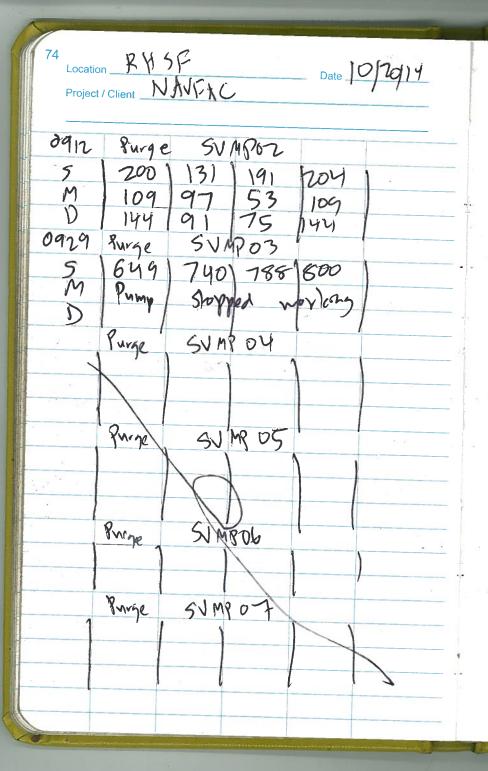
Set up pumpe well.

Set up @ RUMWOZ.

Pump controller still bring.

De pair tunnel will bring.

Other controllers pumping.



Location RHSF

Project / Client NAVFAC

Date 10/21/14

75

Rupole C-W sompling

OBOU C Red 1111 stem meeting.

OBYS @ RIMW 2254-01

Set up to pump.

Controller not discharging

Ing controller not coming

up table. Ing hickan

controller, Dorsn't power on

1200 Leave RIMW 2254-01,

Unable to collect sample.

Vill return Thousday @ 0945,

Depart Red Hill to may

Sittings for pump.

1944 Try to sample RIMDS

but Controller not walnow. 1570 Depart Red 1711 24 DIV' 17

Location PHST- Date 10/27/L 77

Project / Client MVFA C

Personnel 334, KM Calbrate PID

FINTER TUNNED.

Set up a RHMW03

0870 Com DTW- 102.78

7875 Pringed 3 L.

900 Collected ser

d Est. Punged 4 L

Collected sample ES114 ES114790

and ES115 (duje @ 1115).

Set up @ PIJMWO?

DTW= 83.79

Punged 3.5 L

Collected sample ES113.

Depart tunnel.

De 300 1333 1500 1530 1600

Project / Clien NAVPAR

150	-1.01	D			6-2
	7		I CR	188	
S .	13	88	75	00	
D	1.09	126	112	126	95
1013	1 1	09			26
5 1	SUMP	192	105	105	127
M	96	146	122	146	185127
b	163	187	194	194	24
11003	SUMP	04			155
5	197	235	279	279	248
SM	1881	1163	1269	1269	1196
D	85	1105	163	163	129
1124	SUMPO	7	1	1200	
5 M		320	1327	1501	310
Se of the	156	197	1242	242	209
D	1549	522	539	545	
		Ser.		4	
	,				
			1		144
			60	7//	4
			1010	//	
		7/-	1	- /	7 2
				1	

Location R15FProject / Client NAVFACDate 10/28/181

1350 5 M D 1349 5	SVMP 155 13 218 SVMI 1240 1576	193 193 154 18 2031 1992	209 257 278 236 223	200 257 278 278 238 223	
1403 5 1 M D	5VMP 197 143 164	290 290 218 206	323 236 236	323 236 236	253 208 241
	5	A 128	114		
		//			10



APPENDIX C Laboratory Reports





Calscience



WORK ORDER NUMBER: 14-10-2131

The difference is service



AIR | SOIL | WATER | MARINE CHEMISTRY

Analytical Report For

Client: Environmental Science International, Inc.

Client Project Name: Red Hill LTM 112066

Attention: Jeff Hattemer

354 Uluniu Street, Suite 304 Kailua, HI 96734-2500

Richard Vellas

Approved for release on 11/04/2014 by: Richard Villafania

Richard Villatania Project Manager



ResultLink >

Email your PM >

Eurofins Calscience, Inc. (Calscience) certifies that the test results provided in this report meet all NELAC requirements for parameters for which accreditation is required or available. Any exceptions to NELAC requirements are noted in the case narrative. The original report of subcontracted analyses, if any, is attached to this report. The results in this report are limited to the sample(s) tested and any reproduction thereof must be made in its entirety. The client or recipient of this report is specifically prohibited from making material changes to said report and, to the extent that such changes are made, Calscience is not responsible, legally or otherwise. The client or recipient agrees to indemnify Calscience for any defense to any litigation which may arise.



Contents

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Work Order Number:	14-10-2131

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Work Order Narrative

Work Order: 14-10-2131 Page 1 of 1

Condition Upon Receipt:

Samples were received under Chain-of-Custody (COC) on 10/28/14. They were assigned to Work Order 14-10-2131.

Unless otherwise noted on the Sample Receiving forms all samples were received in good condition and within the recommended EPA temperature criteria for the methods noted on the COC. The COC and Sample Receiving Documents are integral elements of the analytical report and are presented at the back of the report.

Holding Times:

All samples were analyzed within prescribed holding times (HT) and/or in accordance with the Calscience Sample Acceptance Policy unless otherwise noted in the analytical report and/or comprehensive case narrative, if required.

Any parameter identified in 40CFR Part 136.3 Table II that is designated as "analyze immediately" with a holding time of <= 15 minutes (40CFR-136.3 Table II, footnote 4), is considered a "field" test and the reported results will be qualified as being received outside of the stated holding time unless received at the laboratory within 15 minutes of the collection time.

Quality Control:

All quality control parameters (QC) were within established control limits except where noted in the QC summary forms or described further within this report.

Additional Comments:

Air - Sorbent-extracted air methods (EPA TO-4A, EPA TO-10, EPA TO-13A, EPA TO-17): Analytical results are converted from mass/sample basis to mass/volume basis using client-supplied air volumes.

New York NELAP air certification does not certify for all reported methods and analytes, reference the accredited items here: http://www.calscience.com/PDF/New_York.pdf

Solid - Unless otherwise indicated, solid sample data is reported on a wet weight basis, not corrected for % moisture. All QC results are always reported on a wet weight basis.

Subcontractor Information:

Unless otherwise noted below (or on the subcontract form), no samples were subcontracted.

Page 1 of 2



Project: Red Hill LTM 112066

Analytical Report

Environmental Science International, Inc. Date Received: 10/28/14 Work Order: 14-10-2131 354 Uluniu Street, Suite 304 Kailua, HI 96734-2500 Preparation: **EPA 3510C** Method: EPA 8015B (M) Units: ug/L

Client Sample Number	Lab Sample Number	Date/Time Collected	Matrix	Instrument	Date Prepared	Date/Time Analyzed	QC Batch ID
ES113	14-10-2131-2-H	10/27/14 14:00	Aqueous	GC 45	10/30/14	10/31/14 09:41	141030B06

Comment(s): - Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are qualified with a "J" flag.

- TPH as Diesel is quantified in the carbon range C10-C28.

<u>DF</u> **Parameter** Result <u>DL</u> <u>LOD</u> **LOQ** Qualifiers TPH as Diesel 12 25 1.00 HD 120 11

Control Limits Qualifiers <u>Surrogate</u> Rec. (%)

n-Octacosane 91 51-141

ES114	14-10-2131-3-H	10/27/14 10:45	Aqueous	GC 45	10/30/14	10/31/14 09:59	141030B06
Comment(s):	- Results were evaluated to the MDL (DL), con	ncentrations >=	to the MDL (D	L) but < RL (LOQ), if found, ar	e qualified with	a "J" flag.
	- TPH as Diesel is quantified in the carbon rar	nge C10-C28.					
<u>Parameter</u>	Result	DL	LOD		LOQ	<u>DF</u>	Qualifiers

Result DL LOD LOQ <u>DF</u> TPH as Diesel 2000 25 1.00 HD 11 12

Surrogate Rec. (%) **Control Limits** Qualifiers

n-Octacosane 92 51-141

ES115	14-10-2131-4-H	10/27/14 11:15	Aqueous	GC 45	10/30/14	10/31/14 10:41	141030B06
Comment(s):	- Results were evaluated to the MDL (DL), con-	centrations >	= to the MDL (DI) but < RL	(LOQ), if found, are	qualified with	a "J" flag.

- TPH as Diesel is quantified in the carbon range C10-C28.

Parameter Result <u>DL</u> LOD LOQ DF Qualifiers TPH as Diesel 2000 11 12 25 1.00 HD

Qualifiers Surrogate **Control Limits** Rec. (%)

n-Octacosane 91 51-141

ES116	14-10-2131-5-H	10/27/14 09:00	Aqueous	GC 45	10/30/14	10/31/14 10:58	141030B06

- Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are qualified with a "J" flag. Comment(s):

- TPH as Diesel is quantified in the carbon range C10-C28.

<u>Parameter</u> LOD LOQ DF Qualifiers Result DL TPH as Diesel 11 12 25 1.00 HD 80

Rec. (%) Surrogate **Control Limits** Qualifiers

65 51-141 n-Octacosane



Environmental Science International, Inc. Date Received: 10/28/14 354 Uluniu Street, Suite 304 Work Order: 14-10-2131 EPA 3510C Kailua, HI 96734-2500 Preparation: Method: EPA 8015B (M) Units: ug/L

Page 2 of 2 Project: Red Hill LTM 112066

Client Sample N	lumber	Lab Sample Number	Date/Time Collected	Matrix	Instrument	Date Prepared	Date/Time Analyzed	QC Batch ID
Method Blank		099-15-516-206	N/A	Aqueous	GC 45	10/30/14	10/31/14 07:03	141030B06
Comment(s):	- Results were evaluated to	the MDL (DL), cond	centrations >= to t	he MDL (DL	_) but < RL (LO	Q), if found, a	re qualified with a	a "J" flag.
<u>Parameter</u>		Result	<u>DL</u>	<u>LOD</u>	LO	<u>2</u>	<u>DF</u>	Qualifiers
TPH as Diesel		<12	11	12	25		1.00	U
Surrogate		Rec. (%)	Control Limi	<u>its Qualit</u>	fiers			
n-Octacosane		117	51-141					





Method Blank

Analytical Report

Environmental Science International, Inc. Date Received: 10/28/14 Work Order: 354 Uluniu Street, Suite 304 14-10-2131 Preparation: EPA 3005A Filt. Kailua, HI 96734-2500 Method: EPA 6020 Units: ug/L

				0				ug,
Project: Red	Hill LTM 112066						P	age 1 of 1
Client Sample N	Number	Lab Sample Number	Date/Time Collected	Matrix	Instrument	Date Prepared	Date/Time Analyzed	QC Batch ID
ES113		14-10-2131-2-G	10/27/14 14:00	Aqueous	ICP/MS 04	10/29/14	10/30/14 23:36	141029L03D
Comment(s):	- Results were evaluated	to the MDL (DL), con	centrations >= t	o the MDL (DL	.) but < RL (LOC), if found, a	are qualified with	a "J" flag.
Parameter Parameter		Result	<u>DL</u>	<u>LOD</u>	LOC	<u>)</u>	<u>DF</u>	Qualifiers
Lead		0.0976	0.0898	0.200	1.00	ı	1.00	J
ES114		14-10-2131-3-G	10/27/14 10:45	Aqueous	ICP/MS 04	10/29/14	10/30/14 23:32	141029L03E
Comment(s):	- Results were evaluated	to the MDL (DL), con	centrations >= t	o the MDL (DL	.) but < RL (LOC	(), if found, a	are qualified with	a "J" flag.
<u>Parameter</u>		Result	<u>DL</u>	<u>LOD</u>	LOC	<u>)</u>	<u>DF</u>	Qualifiers
Lead		<0.200	0.0898	0.200	1.00	l	1.00	U
ES115		14-10-2131-4-G	10/27/14 11:15	Aqueous	ICP/MS 04	10/29/14	10/30/14 23:39	141029L03E
Comment(s):	- Results were evaluated	to the MDL (DL), con	centrations >= t	o the MDL (DL	.) but < RL (LOC	(), if found, a	are qualified with	a "J" flag.
Parameter Parameter		Result	<u>DL</u>	<u>LOD</u>	LOC	<u>)</u>	<u>DF</u>	Qualifiers
Lead		0.165	0.0898	0.200	1.00	l	1.00	J
ES116		14-10-2131-5-G	10/27/14 09:00	Aqueous	ICP/MS 04	10/29/14	10/30/14 23:43	141029L03E
Comment(s):	- Results were evaluated	to the MDL (DL), con	centrations >= t	o the MDL (DL	.) but < RL (LOC)), if found, a	are qualified with	a "J" flag.
Doromotor		Dooult	DI	100	100		DE	Ouglifiers

23110	14-10-2131-3-0	09:00	Aqueous	ICP/WIS 04		23:43	LUSD
Comment(s):	- Results were evaluated to the MDL (DL), con	ncentrations >= t	o the MDL (DL)) but < RL (LOQ)	, if found, are qu	alified with a "J" flag.	
<u>Parameter</u>	Result	<u>DL</u>	LOD	<u>LOQ</u>	<u>DF</u>	Qualifiers	<u> </u>
Lead	<0.200	0.0898	0.200	1.00	1.00) U	

Aqueous

ICP/MS 04

10/29/14

10/30/14

141029L03D

					22.30	
Comment(s):	- Results were evaluated to the MDL (DL	, concentrations >=	to the MDL (DL) but	< RL (LOQ), if fou	nd, are qualified	with a "J" flag.
<u>Parameter</u>	Resul	<u>DL</u>	<u>LOD</u>	<u>LOQ</u>	<u>DF</u>	Qualifiers
Lead	<0.20	0.0898	0.200	1.00	1.00	U

N/A

099-14-497-101



Environmental Science International, Inc. 354 Uluniu Street, Suite 304

Kailua, HI 96734-2500

Date Received: Work Order:

10/28/14 14-10-2131

Preparation:

EPA 3510C

Method:

EPA 8270C SIM PAHs

Units:

ug/L

Project: Red Hill LTM 112066

Page 1 of 5

Client Sample Number	Lab Sample Number	Date/Time Collected	Matrix	Instrument	Date Prepared	Date/Time Analyzed	QC Batch ID
ES113	14-10-2131-2-l	10/27/14 14:00	Aqueous	GC/MS AAA	10/30/14	10/31/14 20:19	141030L08
Comment(s): - Results were evaluated to	to the MDL (DL), con	ncentrations >= to the	he MDL (DL)	but < RL (LOQ), if found, ar	e qualified with a	a "J" flag.
<u>Parameter</u>	Result	<u>DL</u>	LOD	LOQ	<u>)</u>	<u>DF</u>	<u>Qualifiers</u>
Naphthalene	< 0.052	0.035	0.052	0.21		1.00	U
2-Methylnaphthalene	< 0.052	0.048	0.052	0.21		1.00	U
1-Methylnaphthalene	<0.10	0.054	0.10	0.21		1.00	U
Acenaphthylene	< 0.052	0.046	0.052	0.21		1.00	U
Acenaphthene	< 0.052	0.028	0.052	0.21		1.00	U
Fluorene	<0.052	0.044	0.052	0.21		1.00	U
Phenanthrene	<0.052	0.028	0.052	0.21		1.00	U
Anthracene	< 0.052	0.030	0.052	0.21		1.00	U
Fluoranthene	<0.052	0.049	0.052	0.21		1.00	U
Pyrene	< 0.052	0.021	0.052	0.21		1.00	U
Benzo (a) Anthracene	< 0.052	0.034	0.052	0.21		1.00	U
Chrysene	<0.052	0.026	0.052	0.21		1.00	U
Benzo (k) Fluoranthene	<0.052	0.032	0.052	0.21		1.00	U
Benzo (b) Fluoranthene	< 0.052	0.018	0.052	0.21		1.00	U
Benzo (a) Pyrene	<0.052	0.023	0.052	0.21		1.00	U
Indeno (1,2,3-c,d) Pyrene	< 0.052	0.022	0.052	0.21		1.00	U
Dibenz (a,h) Anthracene	< 0.052	0.049	0.052	0.21		1.00	U
Benzo (g,h,i) Perylene	<0.10	0.085	0.10	0.21		1.00	U
<u>Surrogate</u>	Rec. (%)	Control Limi	<u>its Qualifi</u>	<u>ers</u>			
Nitrobenzene-d5	49	28-139					
2-Fluorobiphenyl	65	33-144					
p-Terphenyl-d14	70	23-160					



Environmental Science International, Inc.

Date Received: 10/28/14

354 Uluniu Street, Suite 304

Work Order: 14-10-2131

Kailua, HI 96734-2500

Preparation: EPA 3510C

Preparation: EPA 3510C Method: EPA 8270C SIM PAHs

Units: ug/L
Project: Red Hill LTM 112066 Page 2 of 5

Client Sample Number	Lab Sample Number	Date/Time Collected	Matrix	Instrument	Date Prepared	Date/Time Analyzed	QC Batch ID
ES114	14-10-2131-3-J	10/27/14 10:45	Aqueous	GC/MS AAA	10/30/14	10/31/14 19:10	141030L08
Comment(s): - Results were evaluated	to the MDL (DL), con	centrations >= to t	he MDL (DL	.) but < RL (LOC	Q), if found, a	are qualified with a	a "J" flag.
<u>Parameter</u>	<u>Result</u>	<u>DL</u>	<u>LOD</u>	LOC	<u>3</u>	<u>DF</u>	<u>Qualifiers</u>
Acenaphthylene	<0.047	0.042	0.047	0.19)	1.00	U
Acenaphthene	0.53	0.025	0.047	0.19)	1.00	
Fluorene	< 0.047	0.040	0.047	0.19)	1.00	U
Phenanthrene	<0.047	0.026	0.047	0.19)	1.00	U
Anthracene	< 0.047	0.027	0.047	0.19)	1.00	U
Fluoranthene	<0.047	0.044	0.047	0.19)	1.00	U
Pyrene	< 0.047	0.019	0.047	0.19)	1.00	U
Benzo (a) Anthracene	< 0.047	0.031	0.047	0.19)	1.00	U
Chrysene	<0.047	0.024	0.047	0.19)	1.00	U
Benzo (k) Fluoranthene	<0.047	0.029	0.047	0.19)	1.00	U
Benzo (b) Fluoranthene	< 0.047	0.017	0.047	0.19)	1.00	U
Benzo (a) Pyrene	<0.047	0.021	0.047	0.19)	1.00	U
Indeno (1,2,3-c,d) Pyrene	< 0.047	0.020	0.047	0.19)	1.00	U
Dibenz (a,h) Anthracene	< 0.047	0.045	0.047	0.19)	1.00	U
Benzo (g,h,i) Perylene	<0.095	0.078	0.095	0.19)	1.00	U
<u>Surrogate</u>	Rec. (%)	Control Limi	<u>its Qualif</u>	<u>iers</u>			
Nitrobenzene-d5	66	28-139					
2-Fluorobiphenyl	65	33-144					
p-Terphenyl-d14	68	23-160					

ES114	14-10-2131-3-J	10/27/14 A 10:45	Aqueous	GC/MS AAA	10/30/14	11/03/14 18:28	141030L08
Comment(s): - Results were evaluated to	the MDL (DL), con-	centrations >= to th	e MDL (DL)	but < RL (LOQ)), if found, are	qualified with	a "J" flag.
<u>Parameter</u>	<u>Result</u>	<u>DL</u>	LOD	<u>LOQ</u>	DI	E	Qualifiers
Naphthalene	140	0.64	0.95	3.8	20	0.0	
2-Methylnaphthalene	43	0.88	0.95	3.8	20	0.0	
1-Methylnaphthalene	59	0.98	1.9	3.8	20	0.0	
Surrogate	Rec. (%)	Control Limits	s Qualifie	<u>ers</u>			
Nitrobenzene-d5	60	28-139					
2-Fluorobiphenyl	80	33-144					
p-Terphenyl-d14	83	23-160					



Environmental Science International, Inc. 354 Uluniu Street, Suite 304

Kailua, HI 96734-2500

Date Received: Work Order:

10/28/14 14-10-2131

Preparation:

EPA 3510C

Method:

EPA 8270C SIM PAHs

Units:

ug/L

Project: Red Hill LTM 112066

Page 3 of 5

Client Sample Number	Lab Sample Number	Date/Time Collected	Matrix	Instrument	Date Prepared	Date/Time Analyzed	QC Batch ID
ES115	14-10-2131-4-I	10/27/14 11:15	Aqueous	GC/MS AAA	10/30/14	10/31/14 19:33	141030L08
Comment(s): - Results were evaluated	to the MDL (DL), con	ncentrations >= to t	he MDL (DL	.) but < RL (LOC	(a), if found, a	are qualified with	a "J" flag.
<u>Parameter</u>	<u>Result</u>	<u>DL</u>	LOD	LOC	<u>)</u>	<u>DF</u>	Qualifiers
Acenaphthylene	<0.047	0.042	0.047	0.19)	1.00	U
Acenaphthene	0.53	0.025	0.047	0.19)	1.00	
Fluorene	<0.047	0.040	0.047	0.19)	1.00	U
Phenanthrene	<0.047	0.026	0.047	0.19)	1.00	U
Anthracene	<0.047	0.027	0.047	0.19)	1.00	U
Fluoranthene	<0.047	0.044	0.047	0.19)	1.00	U
Pyrene	<0.047	0.019	0.047	0.19)	1.00	U
Benzo (a) Anthracene	<0.047	0.031	0.047	0.19)	1.00	U
Chrysene	<0.047	0.023	0.047	0.19)	1.00	U
Benzo (k) Fluoranthene	<0.047	0.029	0.047	0.19)	1.00	U
Benzo (b) Fluoranthene	<0.047	0.017	0.047	0.19)	1.00	U
Benzo (a) Pyrene	<0.047	0.021	0.047	0.19)	1.00	U
Indeno (1,2,3-c,d) Pyrene	<0.047	0.020	0.047	0.19)	1.00	U
Dibenz (a,h) Anthracene	<0.047	0.045	0.047	0.19)	1.00	U
Benzo (g,h,i) Perylene	<0.095	0.077	0.095	0.19)	1.00	U
<u>Surrogate</u>	Rec. (%)	Control Limi	its Qualif	<u>iers</u>			
Nitrobenzene-d5	70	28-139					
2-Fluorobiphenyl	69	33-144					
p-Terphenyl-d14	73	23-160					

ES115	14-10-2131-4-I	10/27/14 A 11:15	queous (GC/MS AAA 10/30/1	4 11/03/14 18:52	141030L08
Comment(s): - Results were evaluated to	the MDL (DL), con	centrations >= to the	e MDL (DL)	but < RL (LOQ), if foun	d, are qualified with	a "J" flag.
<u>Parameter</u>	<u>Result</u>	<u>DL</u>	<u>LOD</u>	<u>LOQ</u>	<u>DF</u>	Qualifiers
Naphthalene	130	0.64	0.95	3.8	20.0	
2-Methylnaphthalene	36	0.88	0.95	3.8	20.0	
1-Methylnaphthalene	54	0.98	1.9	3.8	20.0	
<u>Surrogate</u>	Rec. (%)	Control Limits	<u>Qualifie</u>	ers		
Nitrobenzene-d5	53	28-139				
2-Fluorobiphenyl	78	33-144				
p-Terphenyl-d14	80	23-160				



Environmental Science International, Inc. 354 Uluniu Street, Suite 304

Kailua, HI 96734-2500

p-Terphenyl-d14

Date Received: Work Order:

10/28/14 14-10-2131

Preparation:

EPA 3510C EPA 8270C SIM PAHs

Method: Units:

ug/L

Page 4 of 5

Project: Red	Hill LTM	112066
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Client Sample Number	Lab Sample Number	Date/Time Collected	Matrix	Instrument	Date Prepared	Date/Time Analyzed	QC Batch ID
ES116	14-10-2131-5-I	10/27/14 09:00	Aqueous	GC/MS AAA	10/30/14	10/31/14 19:56	141030L08
Comment(s): - Results were evaluated to	o the MDL (DL), con	centrations >= to t	he MDL (DL	.) but < RL (LOC	Q), if found, a	are qualified with	a "J" flag.
<u>Parameter</u>	<u>Result</u>	<u>DL</u>	<u>LOD</u>	LOC	<u>)</u>	<u>DF</u>	<u>Qualifiers</u>
Naphthalene	< 0.054	0.036	0.054	0.21		1.00	U
2-Methylnaphthalene	< 0.054	0.050	0.054	0.21		1.00	U
1-Methylnaphthalene	<0.11	0.055	0.11	0.21		1.00	U
Acenaphthylene	< 0.054	0.048	0.054	0.21		1.00	U
Acenaphthene	< 0.054	0.029	0.054	0.21		1.00	U
Fluorene	< 0.054	0.046	0.054	0.21		1.00	U
Phenanthrene	< 0.054	0.029	0.054	0.21		1.00	U
Anthracene	< 0.054	0.031	0.054	0.21		1.00	U
Fluoranthene	< 0.054	0.050	0.054	0.21		1.00	U
Pyrene	< 0.054	0.022	0.054	0.21		1.00	U
Benzo (a) Anthracene	< 0.054	0.035	0.054	0.21		1.00	U
Chrysene	< 0.054	0.027	0.054	0.21		1.00	U
Benzo (k) Fluoranthene	< 0.054	0.033	0.054	0.21		1.00	U
Benzo (b) Fluoranthene	< 0.054	0.019	0.054	0.21		1.00	U
Benzo (a) Pyrene	< 0.054	0.024	0.054	0.21		1.00	U
Indeno (1,2,3-c,d) Pyrene	< 0.054	0.023	0.054	0.21		1.00	U
Dibenz (a,h) Anthracene	< 0.054	0.051	0.054	0.21		1.00	U
Benzo (g,h,i) Perylene	<0.11	0.088	0.11	0.21	l	1.00	U
Surrogate	Rec. (%)	Control Lim	its Qualif	<u>iers</u>			
Nitrobenzene-d5	59	28-139					
2-Fluorobiphenyl	67	33-144					

23-160



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Kailua, HI 96734-2500

2-Fluorobiphenyl

p-Terphenyl-d14

Date Received: Work Order:

10/28/14 14-10-2131

Preparation:

EPA 3510C

Method:

EPA 8270C SIM PAHs ug/L

Units:

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Project: R	ed Hill	LTM ²	112066
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Client Sample Number	Lab Sample Number	Date/Time Collected	Matrix	Instrument	Date Prepared	Date/Time Analyzed	QC Batch ID
Method Blank	099-15-148-67	N/A	Aqueous	GC/MS AAA	10/30/14	10/31/14 17:17	141030L08
Comment(s): - Results were evaluated to	the MDL (DL), con	centrations >= to the	he MDL (DL)	but < RL (LOC), if found, a	re qualified with a	ı "J" flag.
<u>Parameter</u>	Result	<u>DL</u>	<u>LOD</u>	LOC	<u>)</u>	<u>DF</u>	<u>Qualifiers</u>
Naphthalene	Number Collected Prepared Analyzed		U				
2-Methylnaphthalene	< 0.050	0.046	0.050	0.20		1.00	U
1-Methylnaphthalene	<0.10	0.052	0.10	0.20		1.00	U
Acenaphthylene	< 0.050	0.045	0.050	0.20		1.00	U
Acenaphthene	< 0.050	0.027	0.050	0.20		1.00	U
Fluorene	< 0.050	0.043	0.050	0.20		1.00	U
Phenanthrene	< 0.050	0.027	0.050	0.20		1.00	U
Anthracene	< 0.050	0.029	0.050	0.20		1.00	U
Fluoranthene	< 0.050	0.047	0.050	0.20		1.00	U
Pyrene	< 0.050	0.020	0.050	0.20		1.00	U
Benzo (a) Anthracene	< 0.050	0.033	0.050	0.20		1.00	U
Chrysene	< 0.050	0.025	0.050	0.20		1.00	U
Benzo (k) Fluoranthene	< 0.050	0.031	0.050	0.20		1.00	U
Benzo (b) Fluoranthene	< 0.050	0.018	0.050	0.20		1.00	U
Benzo (a) Pyrene	< 0.050	0.022	0.050	0.20		1.00	U
Indeno (1,2,3-c,d) Pyrene	< 0.050	0.021	0.050	0.20		1.00	U
Dibenz (a,h) Anthracene	< 0.050	0.048	0.050	0.20		1.00	U
Benzo (g,h,i) Perylene	<0.10	0.082	0.10	0.20		1.00	U
Surrogate	Rec. (%)	Control Limi	ts Qualifi	<u>ers</u>			
Nitrobenzene-d5	53	28-139					

33-144

23-160

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Kailua, HI 96734-2500

1,2,4-Trichlorobenzene

Hexachloro-1,3-Butadiene

1,1,1-Trichloroethane

Date Received:

10/28/14 14-10-2131

Work Order: Preparation:

EPA 5030C

Method:

GC/MS / EPA 8260B

Units:

ug/L

Ornio.

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Project: Red Hill LTM 112066

Project: Red Hill LTM 112066							-
Client Sample Number	Lab Sample Number	Date/Time Collected	Matrix	Instrument	Date Prepared	Date/Time Analyzed	QC Batch ID
ES TRIP	14-10-2131-1-A	10/27/14 08:30	Aqueous	GC/MS OO	10/28/14	10/29/14 05:23	141028L046
Comment(s): - Results were evaluated	to the MDL (DL), con	centrations >= t	o the MDL (DI	_) but < RL (LOC	Q), if found, a	re qualified with	a "J" flag.
<u>Parameter</u>	Result	<u>DL</u>	LOD	LOC	<u>3</u>	<u>DF</u>	<u>Qualifiers</u>
Acetone	<10	6.0	10	20		1.00	U,ICH
Benzene	<0.50	0.14	0.50	1.0		1.00	U
Bromodichloromethane	<0.50	0.21	0.50	5.0		1.00	U
Bromoform	<1.0	0.50	1.0	10		1.00	U
Bromomethane	<5.0	3.9	5.0	20		1.00	U
2-Butanone	<5.0	2.2	5.0	10		1.00	U,ICH
Carbon Tetrachloride	<0.50	0.23	0.50	1.0		1.00	U
Chlorobenzene	<0.50	0.17	0.50	5.0		1.00	U
Chloroethane	<5.0	2.3	5.0	10		1.00	U
Chloroform	<0.50	0.46	0.50	5.0		1.00	U
Chloromethane	<2.0	1.8	2.0	10		1.00	U,IJ
Dibromochloromethane	<0.50	0.25	0.50	1.0		1.00	U
1,2-Dibromo-3-Chloropropane	<2.0	1.2	2.0	10		1.00	U
1,2-Dibromoethane	<0.50	0.36	0.50	1.0		1.00	U
1,2-Dichlorobenzene	<0.50	0.46	0.50	1.0		1.00	U
1,3-Dichlorobenzene	<0.50	0.40	0.50	1.0		1.00	U
1,4-Dichlorobenzene	<0.50	0.43	0.50	1.0		1.00	U
1,1-Dichloroethane	<0.50	0.28	0.50	5.0		1.00	U
1,2-Dichloroethane	<0.50	0.24	0.50	1.0		1.00	U,IH
1,1-Dichloroethene	<0.50	0.43	0.50	1.0		1.00	U
c-1,2-Dichloroethene	<0.50	0.48	0.50	1.0		1.00	U
t-1,2-Dichloroethene	<0.50	0.37	0.50	1.0		1.00	U
1,2-Dichloropropane	<0.50	0.42	0.50	5.0		1.00	U
c-1,3-Dichloropropene	<0.50	0.25	0.50	1.0		1.00	U
t-1,3-Dichloropropene	<0.50	0.25	0.50	1.0		1.00	U
Ethylbenzene	<0.50	0.14	0.50	1.0		1.00	U
Methylene Chloride	<1.0	0.64	1.0	5.0		1.00	U
4-Methyl-2-Pentanone	<5.0	4.4	5.0	10		1.00	U
Styrene	<0.50	0.17	0.50	1.0		1.00	U
1,1,1,2-Tetrachloroethane	<0.50	0.40	0.50	1.0		1.00	U
1,1,2,2-Tetrachloroethane	<0.50	0.41	0.50	1.0		1.00	U
Tetrachloroethene	<0.50	0.39	0.50	5.0		1.00	U
Toluene	<0.50	0.24	0.50	1.0		1.00	U

1.0

0.50

0.50

5.0

5.0

1.0

1.00

1.00

1.00

U

U

0.50

0.30

0.32

<1.0

< 0.50

< 0.50



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Date Received: Work Order: Preparation: Method:

Qualifiers

10/28/14 14-10-2131 EPA 5030C GC/MS / EPA 8260B

Units:

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Project: Red Hill LTM 112066

<u>Parameter</u>	Result	<u>DL</u>	<u>LOD</u>	<u>LOQ</u>	<u>DF</u>	Qualifiers
1,1,2-Trichloroethane	<0.50	0.38	0.50	1.0	1.00	U
Trichloroethene	<0.50	0.37	0.50	1.0	1.00	U
1,2,3-Trichloropropane	<1.0	0.64	1.0	5.0	1.00	U
Vinyl Chloride	<0.50	0.30	0.50	1.0	1.00	U
p/m-Xylene	<1.0	0.30	1.0	10	1.00	U
o-Xylene	<0.50	0.23	0.50	1.0	1.00	U
Methyl-t-Butyl Ether (MTBE)	<0.50	0.31	0.50	1.0	1.00	U
Gasoline Range Organics	<30	26	30	50	1.00	U

<u>Surrogate</u>	Rec. (%)	Control Limits
Dibromofluoromethane	99	80-126
1,2-Dichloroethane-d4	101	80-134
Toluene-d8	100	80-120
Toluene-d8-TPPH	99	88-112
1,4-Bromofluorobenzene	94	80-120





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Kailua, HI 96734-2500

Date Received: Work Order:

10/28/14 14-10-2131

Preparation:

EPA 5030C

Method:

GC/MS / EPA 8260B

Units:

ug/L

Project: Red Hill LTM 112066

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Client Sample Number	Lab Sample Number	Date/Time Collected	Matrix	Instrument	Date Date/Tir Prepared Analyze	
ES113	14-10-2131-2-A	10/27/14 14:00	Aqueous	GC/MS OO	10/28/14 10/29/14 06:18	141028L046
Comment(s): - Results were evaluated to	o the MDL (DL), con	centrations >= t	o the MDL (DI	_) but < RL (LOC	Q), if found, are qualified v	vith a "J" flag.
<u>Parameter</u>	Result	<u>DL</u>	LOD	LOC	<u>DF</u>	Qualifiers
Acetone	<10	6.0	10	20	1.00	U,ICH
Benzene	< 0.50	0.14	0.50	1.0	1.00	U
Bromodichloromethane	<0.50	0.21	0.50	5.0	1.00	U
Bromoform	<1.0	0.50	1.0	10	1.00	U
Bromomethane	<5.0	3.9	5.0	20	1.00	U
2-Butanone	<5.0	2.2	5.0	10	1.00	U,ICH
Carbon Tetrachloride	< 0.50	0.23	0.50	1.0	1.00	U
Chlorobenzene	<0.50	0.17	0.50	5.0	1.00	U
Chloroethane	<5.0	2.3	5.0	10	1.00	U
Chloroform	<0.50	0.46	0.50	5.0	1.00	U
Chloromethane	<2.0	1.8	2.0	10	1.00	U,IJ
Dibromochloromethane	<0.50	0.25	0.50	1.0	1.00	U
1,2-Dibromo-3-Chloropropane	<2.0	1.2	2.0	10	1.00	U
1,2-Dibromoethane	<0.50	0.36	0.50	1.0	1.00	U
1,2-Dichlorobenzene	<0.50	0.46	0.50	1.0	1.00	U
1,3-Dichlorobenzene	<0.50	0.40	0.50	1.0	1.00	U
1,4-Dichlorobenzene	<0.50	0.43	0.50	1.0	1.00	U
1,1-Dichloroethane	<0.50	0.28	0.50	5.0	1.00	U
1,2-Dichloroethane	< 0.50	0.24	0.50	1.0	1.00	U,IH
1,1-Dichloroethene	<0.50	0.43	0.50	1.0	1.00	U
c-1,2-Dichloroethene	< 0.50	0.48	0.50	1.0	1.00	U
t-1,2-Dichloroethene	<0.50	0.37	0.50	1.0	1.00	U
1,2-Dichloropropane	<0.50	0.42	0.50	5.0	1.00	U
c-1,3-Dichloropropene	<0.50	0.25	0.50	1.0	1.00	U
t-1,3-Dichloropropene	<0.50	0.25	0.50	1.0	1.00	U
Ethylbenzene	<0.50	0.14	0.50	1.0	1.00	U
Methylene Chloride	<1.0	0.64	1.0	5.0	1.00	U
4-Methyl-2-Pentanone	<5.0	4.4	5.0	10	1.00	U
Styrene	<0.50	0.17	0.50	1.0	1.00	U
1,1,1,2-Tetrachloroethane	<0.50	0.40	0.50	1.0	1.00	U
1,1,2,2-Tetrachloroethane	<0.50	0.41	0.50	1.0	1.00	U
Tetrachloroethene	<0.50	0.39	0.50	5.0	1.00	U
Toluene	<0.50	0.24	0.50	1.0	1.00	U
1,2,4-Trichlorobenzene	<1.0	0.50	1.0	5.0	1.00	U
1,1,1-Trichloroethane	<0.50	0.30	0.50	5.0	1.00	U
Hexachloro-1,3-Butadiene	<0.50	0.32	0.50	1.0	1.00	U



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Work Order: Preparation:

Date Received:

10/28/14 14-10-2131 EPA 5030C

Method:

GC/MS / EPA 8260B

Units:

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Project: Re	ed Hill L	TM 112066
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Toluene-d8

Toluene-d8-TPPH

1,4-Bromofluorobenzene

<u>Parameter</u>	<u>Result</u>	<u>DL</u>	<u>LOD</u>	<u>LOQ</u>	<u>DF</u>	<u>Qualifiers</u>
1,1,2-Trichloroethane	<0.50	0.38	0.50	1.0	1.00	U
Trichloroethene	<0.50	0.37	0.50	1.0	1.00	U
1,2,3-Trichloropropane	<1.0	0.64	1.0	5.0	1.00	U
Vinyl Chloride	<0.50	0.30	0.50	1.0	1.00	U
p/m-Xylene	<1.0	0.30	1.0	10	1.00	U
o-Xylene	<0.50	0.23	0.50	1.0	1.00	U
Methyl-t-Butyl Ether (MTBE)	<0.50	0.31	0.50	1.0	1.00	U
Gasoline Range Organics	<30	26	30	50	1.00	U
Surrogate	Rec. (%)	Control Limits	<u>Qualifiers</u>			
Dibromofluoromethane	95	80-126				
1,2-Dichloroethane-d4	97	80-134				

80-120

88-112

80-120

99

98





Environmental Science International, Inc. 354 Uluniu Street, Suite 304

Kailua, HI 96734-2500

Date Received:

10/28/14 14-10-2131

Work Order: Preparation:

EPA 5030C

Method:

GC/MS / EPA 8260B

Units:

ug/L

Project: Red Hill LTM 112066

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Client Sample Number	Lab Sample Number	Date/Time Collected	Matrix	Instrument	Date Prepared	Date/Time Analyzed	QC Batch ID
ES114	14-10-2131-3-A	10/27/14 10:45	Aqueous	GC/MS OO	10/28/14	10/29/14 05:51	141028L046
Comment(s): - Results were evaluated to	o the MDL (DL), con	centrations >= t	o the MDL (DL	_) but < RL (LOQ), if found, are	qualified with	a "J" flag.
<u>Parameter</u>	Result	<u>DL</u>	LOD	LOQ	<u>D</u>	<u>E</u>	<u>Qualifiers</u>
Acetone	<10	6.0	10	20	1.	00	U,ICH
Benzene	<0.50	0.14	0.50	1.0	1.	00	U
Bromodichloromethane	<0.50	0.21	0.50	5.0	1.	00	U
Bromoform	<1.0	0.50	1.0	10	1.	00	U
Bromomethane	<5.0	3.9	5.0	20	1.	00	U
2-Butanone	<5.0	2.2	5.0	10	1.	00	U,ICH
Carbon Tetrachloride	<0.50	0.23	0.50	1.0	1.	00	U
Chlorobenzene	<0.50	0.17	0.50	5.0	1.	00	U
Chloroethane	<5.0	2.3	5.0	10	1.	00	U
Chloroform	<0.50	0.46	0.50	5.0	1.	00	U
Chloromethane	<2.0	1.8	2.0	10	1.	00	U,IJ
Dibromochloromethane	<0.50	0.25	0.50	1.0	1.	00	U
1,2-Dibromo-3-Chloropropane	<2.0	1.2	2.0	10	1.	00	U
1,2-Dibromoethane	<0.50	0.36	0.50	1.0	1.	00	U
1,2-Dichlorobenzene	<0.50	0.46	0.50	1.0	1.	00	U
1,3-Dichlorobenzene	<0.50	0.40	0.50	1.0	1.	00	U
1,4-Dichlorobenzene	<0.50	0.43	0.50	1.0	1.	00	U
1,1-Dichloroethane	<0.50	0.28	0.50	5.0	1.	00	U
1,2-Dichloroethane	<0.50	0.24	0.50	1.0	1.	00	U,IH
1,1-Dichloroethene	<0.50	0.43	0.50	1.0	1.	00	U
c-1,2-Dichloroethene	<0.50	0.48	0.50	1.0	1.	00	U
t-1,2-Dichloroethene	<0.50	0.37	0.50	1.0	1.	00	U
1,2-Dichloropropane	<0.50	0.42	0.50	5.0	1.	00	U
c-1,3-Dichloropropene	<0.50	0.25	0.50	1.0	1.	00	U
t-1,3-Dichloropropene	<0.50	0.25	0.50	1.0	1.	00	U
Ethylbenzene	0.15	0.14	0.50	1.0	1.	00	J
Methylene Chloride	<1.0	0.64	1.0	5.0	1.	00	U
4-Methyl-2-Pentanone	<5.0	4.4	5.0	10	1.	00	U
Styrene	<0.50	0.17	0.50	1.0	1.	00	U
1,1,1,2-Tetrachloroethane	<0.50	0.40	0.50	1.0	1.	00	U
1,1,2,2-Tetrachloroethane	<0.50	0.41	0.50	1.0	1.	00	U
Tetrachloroethene	<0.50	0.39	0.50	5.0	1.	00	U
Toluene	<0.50	0.24	0.50	1.0	1.	00	U
1,2,4-Trichlorobenzene	<1.0	0.50	1.0	5.0	1.	00	U
1,1,1-Trichloroethane	<0.50	0.30	0.50	5.0	1.	00	U
Hexachloro-1,3-Butadiene	<0.50	0.32	0.50	1.0	1.	00	U



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Date Received: Work Order: Preparation: Method:

10/28/14 14-10-2131 EPA 5030C GC/MS / EPA 8260B

Units:

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ug/L

Project: Red Hill LTM 112066	ò
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1,4-Bromofluorobenzene

<u>Parameter</u>	Result	<u>DL</u>	<u>LOD</u>	<u>LOQ</u>	<u>DF</u>	Qualifiers
1,1,2-Trichloroethane	<0.50	0.38	0.50	1.0	1.00	U
Trichloroethene	<0.50	0.37	0.50	1.0	1.00	U
1,2,3-Trichloropropane	<1.0	0.64	1.0	5.0	1.00	U
Vinyl Chloride	<0.50	0.30	0.50	1.0	1.00	U
p/m-Xylene	<1.0	0.30	1.0	10	1.00	U
o-Xylene	0.32	0.23	0.50	1.0	1.00	J
Methyl-t-Butyl Ether (MTBE)	<0.50	0.31	0.50	1.0	1.00	U
Gasoline Range Organics	57	26	30	50	1.00	
<u>Surrogate</u>	Rec. (%)	Control Limits	<u>Qualifiers</u>			
Dibromofluoromethane	98	80-126				
1,2-Dichloroethane-d4	100	80-134				
Toluene-d8	98	80-120				
Toluene-d8-TPPH	97	88-112				

80-120

10/28/14

14-10-2131



Kailua, HI 96734-2500

Analytical Report

Environmental Science International, Inc.

Date Received:

Work Order:

Preparation: EPA 5030C Method: GC/MS / EPA 8260B

Units: ug/L

Project: Red Hill LTM 112066 Page 7 of 12

Comment(s): - Results were evaluated to the MDL (DL); concentrations >= to the MDL (DL) but < RL (LOQ), if found, are qualified with a *J" flags.	Client Sample Number	Lab Sample Number	Date/Time Collected	Matrix	Instrument	Date Prepared	Date/Time Analyzed	QC Batch ID
Parameter Result DL LOD LOQ DF Qualifiers Acetone <10 6.0 10 20 1.00 UCR Benzene <0.50 0.14 0.50 1.0 1.00 UCR Bromodichloromethane <0.50 0.21 0.50 5.0 1.00 UC Bromoform <1.00 0.50 1.0 10 1.00 UC Bromoform <5.0 3.9 5.0 20 1.00 UC Bromomethane <5.0 2.2 5.0 10 1.00 UJCH Carbon Fetrachloride <5.0 0.23 0.50 1.0 1.00 UJCH Chloroptame <5.0 0.47 0.50 5.0 1.00 U Chloroptame <5.0 0.46 0.50 5.0 1.00 U Chloroptame <2.0 1.2 0.50 1.0 1.00 U Li-Chloroptame <0.50 0.36 <	ES115	14-10-2131-4-A		Aqueous	GC/MS OO	10/28/14		141028L046
Acetone 4.05 6.0 10 20 1.00 U.ICH Benzene 4.05 0.14 0.50 1.0 1.00 U Bromodichioromethane 4.05 0.21 0.50 5.0 1.00 1.00 U Bromomethane 4.50 3.9 5.0 20 1.00 U Ebutanone 4.50 2.2 5.0 10 1.00 U Carbon Tetrachloride 4.050 0.23 0.50 1.0 1.00 U Chlorodorma 4.050 0.17 0.50 5.0 1.00 U Chlorodorma 4.050 0.46 0.50 5.0 1.00 U Chloromethane 4.0 0.46 0.50 5.0 1.00 U Chloromethane 4.0 0.50 0.50 1.0 1.00 U Li-Polibromo-3-Chloropropane 4.20 1.2 2.0 1.0 1.00 U 1,2-Dichlorobenzene 4.050	Comment(s): - Results were evaluated	to the MDL (DL), con	centrations >= t	to the MDL (DI	_) but < RL (LOC	(a), if found, are	qualified with	a "J" flag.
Benzene <0.50 0.14 0.50 1.0 1.00 U Bromodichioromethane <0.50	<u>Parameter</u>	<u>Result</u>	<u>DL</u>	<u>LOD</u>	LOC	<u>D</u>	<u>F</u>	<u>Qualifiers</u>
Bromodichloromethane <0.50 0.21 0.50 5.0 1.00 U Bromoform <1.0	Acetone	<10	6.0	10	20	1	.00	U,ICH
Bromoform <1.0 0.50 1.0 10 1.00 U Bromomethane <5.0	Benzene	< 0.50	0.14	0.50	1.0	1	.00	U
Bromomethane <5.0	Bromodichloromethane	<0.50	0.21	0.50	5.0	1	.00	U
2-Butanone <5.0	Bromoform	<1.0	0.50	1.0	10	1	.00	U
Carbon Tetrachloride <0.50 0.23 0.50 1.0 1.00 U Chlorobenzene <0.50	Bromomethane	<5.0	3.9	5.0	20	1	.00	U
Chlorobenzene <0.50 0.17 0.50 5.0 1.00 U Chlorofethane <5.0	2-Butanone	<5.0	2.2	5.0	10	1	.00	U,ICH
Chloroethane <5.0 2.3 5.0 10 1.00 U Chloroform <0.50	Carbon Tetrachloride	<0.50	0.23	0.50	1.0	1.	.00	U
Chloroform <0.50 0.46 0.50 5.0 1.00 U Chloromethane <2.0	Chlorobenzene	<0.50	0.17	0.50	5.0	1.	.00	U
Chloromethane 42.0 1.8 2.0 10 1.00 U,IJ Dibromochloromethane <0.50	Chloroethane	<5.0	2.3	5.0	10	1	.00	U
Dibromochloromethane <0.50 0.25 0.50 1.0 1.00 U 1,2-Dibromo-3-Chloropropane <2.0	Chloroform	<0.50	0.46	0.50	5.0	1.	.00	U
1,2-Dibromo-3-Chloropropane <2.0	Chloromethane	<2.0	1.8	2.0	10	1.	.00	U,IJ
1,2-Dibromoethane <0.50	Dibromochloromethane	<0.50	0.25	0.50	1.0	1.	.00	U
1,2-Dichlorobenzene <0.50	1,2-Dibromo-3-Chloropropane	<2.0	1.2	2.0	10	1.	.00	U
1,3-Dichlorobenzene <0.50	1,2-Dibromoethane	< 0.50	0.36	0.50	1.0	1.	.00	U
1,4-Dichlorobenzene <0.50	1,2-Dichlorobenzene	<0.50	0.46	0.50	1.0	1.	.00	U
1,1-Dichloroethane <0.50	1,3-Dichlorobenzene	<0.50	0.40	0.50	1.0	1.	.00	U
1,2-Dichloroethane <0.50	1,4-Dichlorobenzene	<0.50	0.43	0.50	1.0	1.	.00	U
1,1-Dichloroethene <0.50	1,1-Dichloroethane	<0.50	0.28	0.50	5.0	1.	.00	U
c-1,2-Dichloroethene <0.50	1,2-Dichloroethane	<0.50	0.24	0.50	1.0	1.	.00	U,IH
t-1,2-Dichloroethene <0.50 0.37 0.50 1.0 1.00 U 1,2-Dichloropropane <0.50	1,1-Dichloroethene	< 0.50	0.43	0.50	1.0	1.	.00	U
1,2-Dichloropropane <0.50	c-1,2-Dichloroethene	<0.50	0.48	0.50	1.0	1.	.00	U
c-1,3-Dichloropropene <0.50	t-1,2-Dichloroethene	<0.50	0.37	0.50	1.0	1.	.00	U
t-1,3-Dichloropropene <0.50	1,2-Dichloropropane	<0.50	0.42	0.50	5.0	1.	.00	U
Ethylbenzene 0.14 0.14 0.50 1.0 1.00 J Methylene Chloride <1.0	c-1,3-Dichloropropene	<0.50	0.25	0.50	1.0	1.	.00	U
Methylene Chloride <1.0 0.64 1.0 5.0 1.00 U 4-Methyl-2-Pentanone <5.0	t-1,3-Dichloropropene	<0.50	0.25	0.50	1.0	1.	.00	U
4-Methyl-2-Pentanone <5.0	Ethylbenzene	0.14	0.14	0.50	1.0	1.	.00	J
Styrene <0.50 0.17 0.50 1.0 1.00 U 1,1,1,2-Tetrachloroethane <0.50	Methylene Chloride	<1.0	0.64	1.0	5.0	1.	.00	U
1,1,1,2-Tetrachloroethane <0.50	4-Methyl-2-Pentanone	<5.0	4.4	5.0	10	1.	.00	U
1,1,2,2-Tetrachloroethane <0.50	Styrene	<0.50	0.17	0.50	1.0	1	.00	U
1,1,2,2-Tetrachloroethane <0.50	1,1,1,2-Tetrachloroethane	<0.50	0.40	0.50	1.0	1	.00	U
Toluene <0.50 0.24 0.50 1.0 1.00 U 1,2,4-Trichlorobenzene <1.0			0.41					
Toluene <0.50 0.24 0.50 1.0 1.00 U 1,2,4-Trichlorobenzene <1.0	Tetrachloroethene	<0.50	0.39	0.50	5.0	1	.00	U
1,2,4-Trichlorobenzene <1.0	Toluene		0.24					U
1,1,1-Trichloroethane <0.50 0.30 0.50 5.0 1.00 U	1,2,4-Trichlorobenzene	<1.0	0.50	1.0	5.0	1	.00	U
	1,1,1-Trichloroethane					1	.00	U
	Hexachloro-1,3-Butadiene			0.50				



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Toluene-d8

Toluene-d8-TPPH

1,4-Bromofluorobenzene

<u>Parameter</u>	<u>Result</u>	<u>DL</u>	<u>LOD</u>	<u>LOQ</u>	<u>DF</u>	Qualifiers
1,1,2-Trichloroethane	<0.50	0.38	0.50	1.0	1.00	U
Trichloroethene	<0.50	0.37	0.50	1.0	1.00	U
1,2,3-Trichloropropane	<1.0	0.64	1.0	5.0	1.00	U
Vinyl Chloride	<0.50	0.30	0.50	1.0	1.00	U
p/m-Xylene	<1.0	0.30	1.0	10	1.00	U
o-Xylene	0.29	0.23	0.50	1.0	1.00	J
Methyl-t-Butyl Ether (MTBE)	<0.50	0.31	0.50	1.0	1.00	U
Gasoline Range Organics	53	26	30	50	1.00	
Surrogate	Rec. (%)	Control Limits	<u>Qualifiers</u>			
Dibromofluoromethane	96	80-126				
1,2-Dichloroethane-d4	97	80-134				

80-120

88-112

80-120

99

98





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Client Sample Number	Lab Sample Number	Date/Time Collected	Matrix	Instrument		Date/Time Q Analyzed	C Batch ID
ES116	14-10-2131-5-A	10/27/14 09:00	Aqueous	GC/MS OO		0/29/14 1/ 07:13	41028L046
Comment(s): - Results were evaluate	d to the MDL (DL), con	centrations >= t	to the MDL (DI	L) but < RL (LOC	Q), if found, are qua	alified with a "J" f	ilag.
<u>Parameter</u>	Result	<u>DL</u>	<u>LOD</u>	LOC	<u>DF</u>	Qu	alifiers
Acetone	<10	6.0	10	20	1.00	U,IC	CH
Benzene	<0.50	0.14	0.50	1.0	1.00	U	
Bromodichloromethane	<0.50	0.21	0.50	5.0	1.00	U	
Bromoform	<1.0	0.50	1.0	10	1.00	U	
Bromomethane	<5.0	3.9	5.0	20	1.00	U	
2-Butanone	<5.0	2.2	5.0	10	1.00	U,IC	CH
Carbon Tetrachloride	<0.50	0.23	0.50	1.0	1.00	U	
Chlorobenzene	<0.50	0.17	0.50	5.0	1.00	U	
Chloroethane	<5.0	2.3	5.0	10	1.00	U	
Chloroform	<0.50	0.46	0.50	5.0	1.00	U	
Chloromethane	<2.0	1.8	2.0	10	1.00	U,I,	J
Dibromochloromethane	<0.50	0.25	0.50	1.0	1.00	U	
1,2-Dibromo-3-Chloropropane	<2.0	1.2	2.0	10	1.00	U	
1,2-Dibromoethane	<0.50	0.36	0.50	1.0	1.00	U	
1,2-Dichlorobenzene	<0.50	0.46	0.50	1.0	1.00	U	
1,3-Dichlorobenzene	<0.50	0.40	0.50	1.0	1.00	U	
1,4-Dichlorobenzene	<0.50	0.43	0.50	1.0	1.00	U	
1,1-Dichloroethane	<0.50	0.28	0.50	5.0	1.00	U	
1,2-Dichloroethane	<0.50	0.24	0.50	1.0	1.00	U,II	Н
1,1-Dichloroethene	<0.50	0.43	0.50	1.0	1.00	U	
c-1,2-Dichloroethene	<0.50	0.48	0.50	1.0	1.00	U	
t-1,2-Dichloroethene	<0.50	0.37	0.50	1.0	1.00	U	
1,2-Dichloropropane	<0.50	0.42	0.50	5.0	1.00	U	
c-1,3-Dichloropropene	<0.50	0.25	0.50	1.0	1.00	U	
t-1,3-Dichloropropene	<0.50	0.25	0.50	1.0	1.00	U	
Ethylbenzene	<0.50	0.14	0.50	1.0	1.00	U	
Methylene Chloride	<1.0	0.64	1.0	5.0	1.00	U	
4-Methyl-2-Pentanone	<5.0	4.4	5.0	10	1.00	U	
Styrene	<0.50	0.17	0.50	1.0	1.00	U	
1,1,1,2-Tetrachloroethane	<0.50	0.40	0.50	1.0	1.00	U	
1,1,2,2-Tetrachloroethane	<0.50	0.41	0.50	1.0	1.00	U	
Tetrachloroethene	<0.50	0.39	0.50	5.0	1.00		
Toluene	<0.50	0.24	0.50	1.0	1.00		
1,2,4-Trichlorobenzene	<1.0	0.50	1.0	5.0	1.00	U	
1,1,1-Trichloroethane	<0.50	0.30	0.50	5.0	1.00	U	
Hexachloro-1,3-Butadiene	<0.50	0.32	0.50	1.0	1.00		



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<u>Parameter</u>	<u>Result</u>	<u>DL</u>	<u>LOD</u>	<u>LOQ</u>	<u>DF</u>	Qualifiers
1,1,2-Trichloroethane	<0.50	0.38	0.50	1.0	1.00	U
Trichloroethene	<0.50	0.37	0.50	1.0	1.00	U
1,2,3-Trichloropropane	<1.0	0.64	1.0	5.0	1.00	U
Vinyl Chloride	<0.50	0.30	0.50	1.0	1.00	U
p/m-Xylene	<1.0	0.30	1.0	10	1.00	U
o-Xylene	<0.50	0.23	0.50	1.0	1.00	U
Methyl-t-Butyl Ether (MTBE)	<0.50	0.31	0.50	1.0	1.00	U
Gasoline Range Organics	<30	26	30	50	1.00	U

Surrogate	Rec. (%)	Control Limits	Qualifiers
Dibromofluoromethane	95	80-126	
1,2-Dichloroethane-d4	98	80-134	
Toluene-d8	99	80-120	
Toluene-d8-TPPH	98	88-112	
1,4-Bromofluorobenzene	95	80-120	





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Client Sample Number	Lab Sample Number	Date/Time Collected	Matrix	Instrument	Date Prepared	Date/Time Analyzed	QC Batch II
Method Blank	099-13-057-71	N/A	Aqueous	GC/MS OO	10/28/14	10/29/14 04:56	141028L04
Comment(s): - Results were evalua	ated to the MDL (DL), cor	ncentrations >=	to the MDL (DI	_) but < RL (LOC	Q), if found, are	e qualified with	a "J" flag.
<u>Parameter</u>	Result	<u>DL</u>	<u>LOD</u>	LOC	<u>)</u>	<u>DF</u>	Qualifiers
Acetone	<10	6.0	10	20	•	1.00	U
Benzene	<0.50	0.14	0.50	1.0	•	1.00	U
Bromodichloromethane	<0.50	0.21	0.50	5.0	•	1.00	U
Bromoform	<1.0	0.50	1.0	10	•	1.00	U
Bromomethane	<5.0	3.9	5.0	20	•	1.00	U
2-Butanone	<5.0	2.2	5.0	10	•	1.00	U
Carbon Tetrachloride	<0.50	0.23	0.50	1.0	•	1.00	U
Chlorobenzene	<0.50	0.17	0.50	5.0		1.00	U
Chloroethane	<5.0	2.3	5.0	10		1.00	U
Chloroform	<0.50	0.46	0.50	5.0		1.00	U
Chloromethane	<2.0	1.8	2.0	10		1.00	U
Dibromochloromethane	<0.50	0.25	0.50	1.0	•	1.00	U
1,2-Dibromo-3-Chloropropane	<2.0	1.2	2.0	10		1.00	U
I,2-Dibromoethane	<0.50	0.36	0.50	1.0		1.00	U
I,2-Dichlorobenzene	<0.50	0.46	0.50	1.0		1.00	U
1,3-Dichlorobenzene	<0.50	0.40	0.50	1.0		1.00	U
1,4-Dichlorobenzene	<0.50	0.43	0.50	1.0		1.00	U
1,1-Dichloroethane	<0.50	0.28	0.50	5.0		1.00	U
1,2-Dichloroethane	<0.50	0.24	0.50	1.0		1.00	U
1,1-Dichloroethene	<0.50	0.43	0.50	1.0		1.00	U
c-1,2-Dichloroethene	<0.50	0.48	0.50	1.0		1.00	U
-1,2-Dichloroethene	<0.50	0.37	0.50	1.0		1.00	U
1,2-Dichloropropane	<0.50	0.42	0.50	5.0		1.00	U
c-1,3-Dichloropropene	<0.50	0.25	0.50	1.0		1.00	U
:-1,3-Dichloropropene	<0.50	0.25	0.50	1.0		1.00	U
Ethylbenzene	<0.50	0.14	0.50	1.0		1.00	U
Methylene Chloride	<1.0	0.64	1.0	5.0		1.00	U
4-Methyl-2-Pentanone	<5.0	4.4	5.0	10		1.00	U
Styrene	<0.50	0.17	0.50	1.0		1.00	U
1,1,1,2-Tetrachloroethane	<0.50	0.40	0.50	1.0		1.00	U
1,1,2,2-Tetrachloroethane	<0.50	0.41	0.50	1.0		1.00	U
Tetrachloroethene	<0.50	0.39	0.50	5.0		1.00	U
Toluene	<0.50	0.24	0.50	1.0		1.00	U
1,2,4-Trichlorobenzene	<1.0	0.50	1.0	5.0		1.00	U
1,1,1-Trichloroethane	<0.50	0.30	0.50	5.0		1.00	U
Hexachloro-1,3-Butadiene	<0.50	0.32	0.50	1.0	•	1.00	U



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Toluene-d8-TPPH

1,4-Bromofluorobenzene

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<u>Parameter</u>	Result	<u>DL</u>	<u>LOD</u>	<u>LOQ</u>	<u>DF</u>	Qualifiers
1,1,2-Trichloroethane	<0.50	0.38	0.50	1.0	1.00	U
Trichloroethene	< 0.50	0.37	0.50	1.0	1.00	U
1,2,3-Trichloropropane	<1.0	0.64	1.0	5.0	1.00	U
Vinyl Chloride	<0.50	0.30	0.50	1.0	1.00	U
p/m-Xylene	<1.0	0.30	1.0	10	1.00	U
o-Xylene	< 0.50	0.23	0.50	1.0	1.00	U
Methyl-t-Butyl Ether (MTBE)	< 0.50	0.31	0.50	1.0	1.00	U
Gasoline Range Organics	<30	26	30	50	1.00	U
<u>Surrogate</u>	Rec. (%)	Control Limits	<u>Qualifiers</u>			
Dibromofluoromethane	98	80-126				
1,2-Dichloroethane-d4	101	80-134				
Toluene-d8	100	80-120				

88-112

80-120

100



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Preparation:

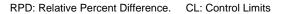
EPA 3510C

Method:

EPA 8015B (M)

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Quality Control Sample ID	Туре		Matrix	In	strument	Date Prepared	Date Ana	lyzed	MS/MSD Ba	tch Number
ES114	Sample	Sample		Aqueous GC 45		10/30/14	10/31/14	09:59	141030S06	
ES114	Matrix Spike	Matrix Spike Aqu		Aqueous GC 45		10/30/14	14 10/31/14 09:05		141030S06	
ES114	Matrix Spike	latrix Spike Duplicate		Aqueous GC 45		10/30/14	10/31/14	09:23	141030S06	
Parameter	<u>Sample</u> <u>Conc.</u>	<u>Spike</u> Added	MS Conc.	MS %Rec.	MSD Conc.	MSD %Rec.	%Rec. CL	RPD	RPD CL	Qualifiers
TPH as Diesel	2046	2000	7915	293	7994	297	55-133	1	0-30	3





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Preparation:

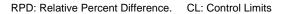
Method:

EPA 3005A Filt.

Method:

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Quality Control Sample ID	Туре		Matrix	Ir	nstrument	Date Prepared	Date Ana	lyzed	MS/MSD Bat	ch Number
ES114	Sample		Aqueou	s IC	CP/MS 04	10/29/14	10/30/14	23:32	141029S03B	
ES114	Matrix Spike		Aqueou	s IC	CP/MS 04	10/29/14	10/30/14	23:05	141029S03B	
ES114	Matrix Spike	Duplicate	Aqueou	s IC	CP/MS 04	10/29/14	10/30/14	23:19	141029S03B	
Parameter	Sample Conc.	<u>Spike</u> Added	MS Conc.	MS %Rec.	MSD Conc.	MSD %Rec.	%Rec. CL	RPD	RPD CL	Qualifiers
Lead	ND	100.0	112.2	112	112.7	113	80-120	0	0-20	





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EPA 3510C

EPA 8270C SIM PAHs

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Quality Control Sample ID	Туре		Matrix	l	nstrument	Date Prepa	red Date Ana	lyzed	MS/MSD Ba	tch Number
ES114	Sample		Aqueou	ıs (GC/MS AAA	10/30/14	11/03/14	18:28	141030S08	
ES114	Matrix Spike		Aqueou	ıs (GC/MS AAA	10/30/14	10/31/14	18:04	141030S08	
ES114	Matrix Spike	Duplicate	Aqueou	ıs (GC/MS AAA	10/30/14	10/31/14	18:27	141030 S 08	
Parameter	Sample Conc.	<u>Spike</u> <u>Added</u>	MS Conc.	MS %Rec	MSD Conc.	MSD %Rec.	%Rec. CL	RPD	RPD CL	Qualifiers
Naphthalene	143.3	2.000	83.89	0	78.94	0	21-133	6	0-25	3
2-Methylnaphthalene	42.74	2.000	34.71	0	32.01	0	21-140	8	0-25	3
1-Methylnaphthalene	59.00	2.000	47.61	0	45.00	0	20-140	6	0-25	3
Acenaphthylene	ND	2.000	1.706	85	1.549	77	33-145	10	0-25	
Acenaphthene	0.5300	2.000	2.109	79	1.970	72	49-121	7	0-25	
Fluorene	ND	2.000	1.860	93	1.701	85	59-121	9	0-25	
Phenanthrene	ND	2.000	1.624	81	1.459	73	54-120	11	0-25	
Anthracene	ND	2.000	1.548	77	1.417	71	27-133	9	0-25	
Fluoranthene	ND	2.000	1.666	83	1.524	76	26-137	9	0-25	
Pyrene	ND	2.000	1.581	79	1.473	74	18-168	7	0-25	
Benzo (a) Anthracene	ND	2.000	1.551	78	1.422	71	33-143	9	0-25	
Chrysene	ND	2.000	1.540	77	1.402	70	17-168	9	0-25	
Benzo (k) Fluoranthene	ND	2.000	1.447	72	1.319	66	24-159	9	0-25	
Benzo (b) Fluoranthene	ND	2.000	1.372	69	1.248	62	24-159	9	0-25	
Benzo (a) Pyrene	ND	2.000	1.331	67	1.216	61	17-163	9	0-25	
Indeno (1,2,3-c,d) Pyrene	ND	2.000	1.320	66	1.166	58	10-171	12	0-25	
Dibenz (a,h) Anthracene	ND	2.000	1.422	71	1.275	64	10-219	11	0-25	
Benzo (g,h,i) Perylene	ND	2.000	1.524	76	1.373	69	10-227	10	0-25	



Environmental Science International, Inc.

Date Received:

Work Order:

10/28/14

Work Order:

14-10-2131

Kailua, HI 96734-2500

Preparation:

EPA 5030C

Method: GC/MS / EPA 8260B

Project: Red Hill LTM 112066	Page 4 of 5
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Quality Control Sample ID	Туре		Matrix	Ir	nstrument	Date Prepared	Date Ana	lyzed	MS/MSD Ba	tch Number
ES114	Sample		Aqueous	G	C/MS OO	10/28/14	10/29/14	05:51	141028S037	
ES114	Matrix Spike		Aqueous	G	C/MS OO	10/28/14	10/29/14	07:40	141028S037	
ES114	Matrix Spike	Duplicate	Aqueous	G	C/MS OO	10/28/14	10/29/14	08:07	141028S037	
<u>Parameter</u>	Sample Conc.	Spike Added	MS Conc.	MS %Rec.	MSD Conc.	MSD %Rec.	%Rec. CL	RPD	RPD CL	Qualifiers
Acetone	ND	50.00	55.96	112	59.02	118	40-140	5	0-20	
Benzene	ND	50.00	48.83	98	47.07	94	80-120	4	0-20	
Bromodichloromethane	ND	50.00	48.14	96	47.21	94	75-120	2	0-20	
Bromoform	ND	50.00	46.72	93	46.10	92	70-130	1	0-20	
Bromomethane	ND	50.00	47.90	96	42.68	85	30-145	12	0-20	
2-Butanone	ND	50.00	49.72	99	54.71	109	30-150	10	0-20	
Carbon Tetrachloride	ND	50.00	41.49	83	41.52	83	65-140	0	0-20	
Chlorobenzene	ND	50.00	49.02	98	47.13	94	80-120	4	0-20	
Chloroethane	ND	50.00	48.97	98	46.89	94	60-135	4	0-20	
Chloroform	ND	50.00	48.66	97	47.25	94	65-135	3	0-20	
Chloromethane	ND	50.00	42.58	85	42.98	86	40-125	1	0-20	
Dibromochloromethane	ND	50.00	49.94	100	48.76	98	60-135	2	0-20	
1,2-Dibromo-3-Chloropropane	ND	50.00	42.13	84	42.15	84	50-130	0	0-20	
1,2-Dibromoethane	ND	50.00	50.33	101	49.09	98	80-120	2	0-20	
1,2-Dichlorobenzene	ND	50.00	47.77	96	46.40	93	70-120	3	0-20	
1,3-Dichlorobenzene	ND	50.00	47.15	94	45.84	92	75-125	3	0-20	
1,4-Dichlorobenzene	ND	50.00	46.75	94	45.19	90	75-125	3	0-20	
1,1-Dichloroethane	ND	50.00	45.35	91	44.62	89	70-135	2	0-20	
1,2-Dichloroethane	ND	50.00	54.76	110	52.87	106	70-130	4	0-20	
1,1-Dichloroethene	ND	50.00	46.83	94	46.17	92	70-130	1	0-20	
c-1,2-Dichloroethene	ND	50.00	48.30	97	47.53	95	70-125	2	0-20	
t-1,2-Dichloroethene	ND	50.00	46.93	94	46.11	92	60-140	2	0-20	
1,2-Dichloropropane	ND	50.00	47.05	94	46.05	92	75-125	2	0-20	
c-1,3-Dichloropropene	ND	50.00	43.96	88	43.95	88	70-130	0	0-20	
t-1,3-Dichloropropene	ND	50.00	37.58	75	38.14	76	55-140	1	0-20	
Ethylbenzene	ND	50.00	48.57	97	46.29	93	75-125	5	0-20	
Methylene Chloride	ND	50.00	46.74	93	46.27	93	55-140	1	0-20	
4-Methyl-2-Pentanone	ND	50.00	45.66	91	46.54	93	60-135	2	0-20	
Styrene	ND	50.00	49.11	98	47.51	95	65-135	3	0-20	
1,1,1,2-Tetrachloroethane	ND	50.00	46.10	92	45.50	91	80-130	1	0-20	
1,1,2,2-Tetrachloroethane	ND	50.00	48.52	97	47.00	94	65-130	3	0-20	
Tetrachloroethene	ND	50.00	53.01	106	50.86	102	45-150	4	0-20	
Toluene	ND	50.00	48.79	98	47.39	95	75-120	3	0-20	
1,2,4-Trichlorobenzene	ND	50.00	47.10	94	46.48	93	65-135	1	0-20	
1,1,1-Trichloroethane	ND	50.00	43.81	88	43.80	88	65-130	0	0-20	



Environmental Science International, Inc.

Date Received:

Work Order:

14-10-2131

Kailua, HI 96734-2500

Preparation:

Method:

Date Received:

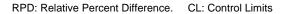
10/28/14

Preparation:

EPA 5030C

Project: Red Hill LTM 112066 Page 5 of 5

<u>Parameter</u>	Sample Conc.	Spike Added	MS Conc.	MS %Rec.	MSD Conc.	MSD %Rec.	%Rec. CL	RPD	RPD CL	Qualifiers
Hexachloro-1,3-Butadiene	ND	50.00	47.57	95	45.21	90	50-140	5	0-20	
1,1,2-Trichloroethane	ND	50.00	47.82	96	46.98	94	75-125	2	0-20	
Trichloroethene	ND	50.00	47.24	94	44.84	90	70-125	5	0-20	
1,2,3-Trichloropropane	ND	50.00	50.56	101	49.71	99	75-125	2	0-20	
Vinyl Chloride	ND	50.00	51.75	103	51.09	102	50-145	1	0-20	
p/m-Xylene	ND	100.0	96.07	96	92.12	92	75-130	4	0-20	
o-Xylene	ND	50.00	49.17	98	47.56	95	80-120	3	0-20	
Methyl-t-Butyl Ether (MTBE)	ND	50.00	40.87	82	42.14	84	65-125	3	0-20	





Quality Control - PDS

Environmental Science International, Inc.

Date Received:

Work Order:

10/28/14

Work Order:

14-10-2131

Kailua, HI 96734-2500

Preparation:

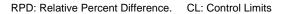
Method:

EPA 3005A Filt.

Method:

Project: Red Hill LTM 112066 Page 1 of 1

Quality Control Sample ID	Туре	N	/latrix	Instrument	Date Prepared	Date Analyzed	PDS/PDSD Batch Number
ES114	Sample	Δ.	Aqueous	ICP/MS 04	10/29/14 00:00	10/30/14 23:32	141029S03B
ES114	PDS	A	Aqueous	ICP/MS 04	10/29/14 00:00	10/30/14 23:22	141029S03B
<u>Parameter</u>		Sample Conc.	Spike Added	PDS Conc.	PDS %Re	ec. %Rec. C	<u>CL</u> <u>Qualifiers</u>
Lead		ND	100.0	106.6	107	75-125	





Quality Control - LCS/LCSD

Environmental Science International, Inc.

Date Received:

Work Order:

10/28/14

Work Order:

14-10-2131

Kailua, HI 96734-2500

Preparation:

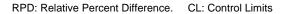
EPA 3510C

Method:

EPA 8015B (M)

Project: Red Hill LTM 112066 Page 1 of 5

Quality Control Sample ID	Туре	Mat	rix	Instrument	Date Prep	pared Date	Analyzed	LCS/LCSD Ba	atch Number
099-15-516-206	LCS	Aqı	ieous	GC 45	10/30/14	10/3	1/14 07:20	141030B06	
099-15-516-206	LCSD	Aqu	ueous	GC 45	10/30/14	10/3	1/14 07:37	141030B06	
Parameter	Spike Added	LCS Conc.	LCS %Rec.	LCSD Conc.	LCSD %Rec.	%Rec. CL	RPD	RPD CL	Qualifiers
TPH as Diesel	2000	2296	115	2219	111	60-132	3	0-11	





Quality Control - LCS

Environmental Science International, Inc.

Date Received:

Work Order:

10/28/14

Work Order:

14-10-2131

Kailua, HI 96734-2500

Preparation:

Method:

EPA 3005A Filt.

Method:

Project: Red Hill LTM 112066 Page 2 of 5

Quality Control Sample ID	Туре	Matrix	Instrument	Date Prepared	Date Analyzed	LCS Batch Number
099-14-497-101	LCS	Aqueous	ICP/MS 04	10/29/14	10/30/14 23:02	141029L03D
Parameter		Spike Added	Conc. Recover	ed LCS %Re	ec. %Rec	. CL Qualifiers
Lead		100.0	106.3	106	80-120)



Quality Control - LCS

Environmental Science International, Inc. 354 Uluniu Street, Suite 304

Kailua, HI 96734-2500

Project: Red Hill LTM 112066

Date Received:

Method:

Work Order: 14-10-2131 EPA 3510C

Preparation:

EPA 8270C SIM PAHs

10/28/14

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Quality Control Sample ID	Туре	Matrix	Instrument	Date Prepared	Date Analyzed LC	S Batch Number
099-15-148-67	LCS	Aqueous	GC/MS AAA	10/30/14	10/31/14 17:40 141	030L08
<u>Parameter</u>		Spike Added	Conc. Recov	ered LCS %Re	ec. %Rec. CL	Qualifiers
Naphthalene		2.000	1.394	70	21-133	
2-Methylnaphthalene		2.000	1.382	69	21-140	
1-Methylnaphthalene		2.000	1.409	70	20-140	
Acenaphthylene		2.000	1.394	70	33-145	
Acenaphthene		2.000	1.451	73	55-121	
Fluorene		2.000	1.539	77	59-121	
Phenanthrene		2.000	1.447	72	54-120	
Anthracene		2.000	1.308	65	27-133	
Fluoranthene		2.000	1.538	77	26-137	
Pyrene		2.000	1.480	74	45-129	
Benzo (a) Anthracene		2.000	1.398	70	33-143	
Chrysene		2.000	1.425	71	17-168	
Benzo (k) Fluoranthene		2.000	1.412	71	24-159	
Benzo (b) Fluoranthene		2.000	1.322	66	24-159	
Benzo (a) Pyrene		2.000	1.150	58	17-163	
Indeno (1,2,3-c,d) Pyrene		2.000	1.251	63	25-175	
Dibenz (a,h) Anthracene		2.000	1.391	70	25-175	
Benzo (g,h,i) Perylene		2.000	1.514	76	25-157	

RPD: Relative Percent Difference. CL: Control Limits



Quality Control - LCS/LCSD

Environmental Science International, Inc. 354 Uluniu Street, Suite 304

Kailua, HI 96734-2500

Project: Red Hill LTM 112066

Date Received: Work Order:

14-10-2131

10/28/14

Preparation:

EPA 5030C GC/MS / EPA 8260B

Method: GC/MS / EPA 8260 Page 4 of 5

Quality Control Sample ID	Туре	Ma	trix	Instrument	Date Pre	pared I	Date Analyze	ed_LCS/LCSD Ba	atch Number
099-13-057-71	LCS	Aq	ueous	GC/MS OO	10/28/14		10/29/14 03:	34 141028L046	
099-13-057-71	LCSD	Aq	ueous	GC/MS OO	10/28/14		10/29/14 04:	01 141028L046	
Parameter	Spike Added	LCS Conc.	LCS %Rec.	LCSD Conc.	LCSD %Rec.	%Rec.	CL RPD	RPD CL	Qualifiers
Acetone	50.00	44.78	90	N/A	N/A	40-140	N/A	0-20	
Benzene	50.00	51.34	103	N/A	N/A	80-120	N/A	0-20	
Bromodichloromethane	50.00	51.61	103	N/A	N/A	75-120	N/A	0-20	
Bromoform	50.00	49.35	99	N/A	N/A	70-130	N/A	0-20	
Bromomethane	50.00	44.25	88	N/A	N/A	30-145	N/A	0-20	
2-Butanone	50.00	47.46	95	N/A	N/A	30-150	N/A	0-20	
Carbon Tetrachloride	50.00	45.03	90	N/A	N/A	65-140	N/A	0-20	
Chlorobenzene	50.00	50.81	102	N/A	N/A	80-120	N/A	0-20	
Chloroethane	50.00	52.54	105	N/A	N/A	60-135	N/A	0-20	
Chloroform	50.00	51.69	103	N/A	N/A	65-135	N/A	0-20	
Chloromethane	50.00	44.16	88	N/A	N/A	40-125	N/A	0-20	
Dibromochloromethane	50.00	52.37	105	N/A	N/A	60-135	N/A	0-20	
1,2-Dibromo-3-Chloropropane	50.00	41.53	83	N/A	N/A	50-130	N/A	0-20	
1,2-Dibromoethane	50.00	52.74	105	N/A	N/A	80-120	N/A	0-20	
1,2-Dichlorobenzene	50.00	49.45	99	N/A	N/A	70-120	N/A	0-20	
1,3-Dichlorobenzene	50.00	48.61	97	N/A	N/A	75-125	N/A	0-20	
1,4-Dichlorobenzene	50.00	48.44	97	N/A	N/A	75-125	N/A	0-20	
1,1-Dichloroethane	50.00	48.98	98	N/A	N/A	70-135	N/A	0-20	
1,2-Dichloroethane	50.00	58.50	117	N/A	N/A	70-130	N/A	0-20	
1,1-Dichloroethene	50.00	50.88	102	N/A	N/A	70-130	N/A	0-20	
c-1,2-Dichloroethene	50.00	51.67	103	N/A	N/A	70-125	N/A	0-20	
t-1,2-Dichloroethene	50.00	50.30	101	N/A	N/A	60-140	N/A	0-20	
1,2-Dichloropropane	50.00	50.03	100	N/A	N/A	75-125	N/A	0-20	
c-1,3-Dichloropropene	50.00	48.68	97	N/A	N/A	70-130	N/A	0-20	
t-1,3-Dichloropropene	50.00	43.01	86	N/A	N/A	55-140	N/A	0-20	
Ethylbenzene	50.00	50.41	101	N/A	N/A	75-125	N/A	0-20	
Methylene Chloride	50.00	50.23	100	N/A	N/A	55-140	N/A	0-20	
4-Methyl-2-Pentanone	50.00	49.12	98	N/A	N/A	60-135	N/A	0-20	
Styrene	50.00	50.88	102	N/A	N/A	65-135	N/A	0-20	
1,1,1,2-Tetrachloroethane	50.00	48.61	97	N/A	N/A	80-130	N/A	0-20	
1,1,2,2-Tetrachloroethane	50.00	49.80	100	N/A	N/A	65-130	N/A	0-20	
Tetrachloroethene	50.00	52.93	106	N/A	N/A	45-150	N/A	0-20	
Toluene	50.00	51.72	103	N/A	N/A	75-120	N/A	0-20	
1,2,4-Trichlorobenzene	50.00	48.11	96	N/A	N/A	65-135	N/A	0-20	
1,1,1-Trichloroethane	50.00	47.64	95	N/A	N/A	65-130	N/A	0-20	
Hexachloro-1,3-Butadiene	50.00	48.39	97	N/A	N/A	50-140	N/A	0-20	



Quality Control - LCS/LCSD

Environmental Science International, Inc.

Date Received:

Work Order:

10/28/14

Work Order:

14-10-2131

Kailua, HI 96734-2500

Preparation:

Method:

GC/MS / EPA 8260B

Project: Red Hill LTM 112066 Page 5 of 5

<u>Parameter</u>	Spike Added	LCS Conc.	LCS %Rec.	LCSD Conc.	LCSD %Rec.	%Rec. CL	RPD	RPD CL	Qualifiers
1,1,2-Trichloroethane	50.00	50.32	101	N/A	N/A	75-125	N/A	0-20	
Trichloroethene	50.00	49.50	99	N/A	N/A	70-125	N/A	0-20	
1,2,3-Trichloropropane	50.00	52.64	105	N/A	N/A	75-125	N/A	0-20	
Vinyl Chloride	50.00	54.73	109	N/A	N/A	50-145	N/A	0-20	
p/m-Xylene	100.0	100.3	100	N/A	N/A	75-130	N/A	0-20	
o-Xylene	50.00	51.10	102	N/A	N/A	80-120	N/A	0-20	
Methyl-t-Butyl Ether (MTBE)	50.00	45.38	91	N/A	N/A	65-125	N/A	0-20	
Gasoline Range Organics	1000	1077	108	1032	103	80-120	4	0-20	



Sample Analysis Summary Report

Work Order: 14-10-2131				Page 1 of 1
Method	Extraction	Chemist ID	Instrument	Analytical Location
EPA 6020	EPA 3005A Filt.	598	ICP/MS 04	1
EPA 8015B (M)	EPA 3510C	682	GC 45	1
EPA 8270C SIM PAHs	EPA 3510C	923	GC/MS AAA	1
GC/MS / EPA 8260B	EPA 5030C	849	GC/MS OO	2

Location 1: 7440 Lincoln Way, Garden Grove, CA 92841 Location 2: 7445 Lampson Avenue, Garden Grove, CA 92841



Glossary of Terms and Qualifiers

Work Order: 14-10-2131 Page 1 of 1

Qualifiers	<u>Definition</u>
*	See applicable analysis comment.
<	Less than the indicated value.
>	Greater than the indicated value.
1	Surrogate compound recovery was out of control due to a required sample dilution. Therefore, the sample data was reported without further clarification.
2	Surrogate compound recovery was out of control due to matrix interference. The associated method blank surrogate spike compound was in control and, therefore, the sample data was reported without further clarification.
3	Recovery of the Matrix Spike (MS) or Matrix Spike Duplicate (MSD) compound was out of control due to suspected matrix interference. The associated LCS recovery was in control.
4	The MS/MSD RPD was out of control due to suspected matrix interference.
5	The PDS/PDSD or PES/PESD associated with this batch of samples was out of control due to suspected matrix interference.
6	Surrogate recovery below the acceptance limit.
7	Surrogate recovery above the acceptance limit.
В	Analyte was present in the associated method blank.
BU	Sample analyzed after holding time expired.
BV	Sample received after holding time expired.
DL	The Detection Limit (DL) is the smallest analyte concentration that can be demonstrated to be different from zero or a blank concentration at the 99% level of confidence.
Е	Concentration exceeds the calibration range.
ET	Sample was extracted past end of recommended max. holding time.
HD	The chromatographic pattern was inconsistent with the profile of the reference fuel standard.
ICH	Initial calibration verification recovery is above the control limit for this analyte.
ICJ	Initial calibration verification recovery is below the control limit for this analyte.
IH	Calibration verification recovery is above the control limit for this analyte.
IJ	Calibration verification recovery is below the control limit for this analyte.
J	Analyte was detected at a concentration below the LOQ and above the DL. Reported value is estimated.
JA	Analyte positively identified but quantitation is an estimate.
LOD	The Limit of Detection (LOD) is the smallest amount or concentration of a substance that must be present in a sample in order to be detected at 99% confidence level.
LOQ	The Limit of Quantitation (LOQ) is the lowest concentration of a substance that produces a quantitative result within specified limits of precision and bias.
Q	Spike recovery and RPD control limits do not apply resulting from the parameter concentration in the sample exceeding the spike concentration by a factor of four or greater.

- SG The sample extract was subjected to Silica Gel treatment prior to analysis.
- Undetected at Detection Limit (DL) and is reported as less than the Limit of Detection (LOD).
- X % Recovery and/or RPD out-of-range.
- Z Analyte presence was not confirmed by second column or GC/MS analysis.

Solid - Unless otherwise indicated, solid sample data is reported on a wet weight basis, not corrected for % moisture. All QC results are reported on a wet weight basis.

Any parameter identified in 40CFR Part 136.3 Table II that is designated as "analyze immediately" with a holding time of <= 15 minutes (40CFR-136.3 Table II, footnote 4), is considered a "field" test and the reported results will be qualified as being received outside of the stated holding time unless received at the laboratory within 15 minutes of the collection time.

A calculated total result (Example: Total Pesticides) is the summation of each component concentration and/or, if "J" flags are reported, estimated concentration. Component concentrations showing not detected (ND) are summed into the calculated total result as zero concentrations.

alscience 7440 Lincol Other location CLIENT: ESI MA 161-67-40 TIME (Rush surcharge	Ilence Environmental Laboratories, Inc. 7440 Lincoln Way, Garden Grove, CA 92841-1427 • (714) 895-5494 Other locations: Concord, San Luis Obispo, Houston, and Corpus Christi For courier service / sample drop off information, contact sales@calscience.com or call us. 146 140 140 140 140 140 140 140	841-1427 9, Houstor Do off info COM or ca H S' H S' H S' H S' WDARD"):	27 • (714) con, and Co information, call us. STATE:	1101 895-54 orpus C	1895-5494 Orpus Christi		WO#/L/	WO#/LAB USE ONLY LA-10-213 CLIENT PROJECT NAME/N PROJECT CONTACT: "Pe Jeff H Please chec	SE ONLY ECT NAM I U NTACT: XAF		NUMBER: 1 120 L Hightemer REQUES	Date Date INMBER: 120 CC 20 CC 20 CC 3 CC 4 CC 5 CC 6 CC 7 CC	3 3 3	NAL S	ANALYSES	OF CUSTON OF P.O. NO.: VSES OG.	B Sunt	ANALYSES TAIN OF CUSTODY RECORD Of Of Of Other Control	
SAME DAY 24 HR 18 HR 172 HR 15 DAYS X GLOBAL ID SPECIAL INSTRUCTIONS: Regular (full) (1M Minitoring analyte 1959) Regular (full) (1M minitoring analyte 1959)	172 HR DE	Mayte i	X Z	STANDARD L NO. NO. OF	Ibueseuved 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0	sserved A Filtered	(0)78) 089. (6)Hd1	н Се-сзе Се-с 44		EX / MTBE = 8260 = ==================================	ygenates (8260)	sp (5035) □ En Core □ Terra Core	OCs (8270)	Bs (8082)	MI2 0728 ½ 0728 □ 2H	XT4T\0508 \(\text{XT4T\0108} \(\text{C020\0174} \)	0100 6170 61		
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DISTRIBUTION: White with final report, Green and Yellow to Client.

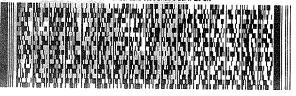
Please note that pages 1 and 2 of 2 of our T/Cs are printed on the reverse side of the Green and Yellow copies, respectively.

Page 37 of 40



SHIP DATE: 270CT14 ACTWGT: 56.1 LB CAD: /POS1525 DIMS: 26x14x13 IN

BILL RECIPIENT



FedEx Express

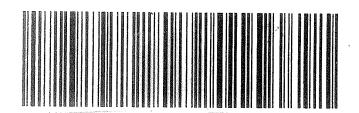


1 of 2 TRK# 0200 8045 5791 7364

TUE - 28 OCT AA STANDARD OVERNIGHT

WZ APVA

92841 CA-US SNA



SHIP DATE: 270CT14 ACTWGT: 48.4 LB CAD: /POS1525 DIMS: 26x14x13 IN

UNITED STATES US

SAMPLE CONTROL CALSCIENCE 7440 LINCOLN WAY

GARDEN GROVE CA 92841

FedEx Express

2 of 2 MPS# 7801 5009 0734 Mstr# 8045 5791 7364

TUE - 28 OCT AA STANDARD OVERNIGHT

0200

WZ APVA

92841

SNA





Calscience

WORK ORDER #: 14-10- 2 1 3 1

SAMPLE RECEIPT FORM

Cooler ___ of ___

CLIENT:EST	DATE: _	10/28/	14
TEMPERATURE: Thermometer ID: SC2 (Criteria: 0.0 °C – 6.0 °C, not froze	n except se	diment/tissue)	
	Blank	☐ Sample	
☐ Sample(s) outside temperature criteria (PM/APM contacted by:)			
☐ Sample(s) outside temperature criteria but received on ice/chilled on same of	lay of sampl	ing.	
☐ Received at ambient temperature, placed on ice for transport by Co		Ü	
Ambient Temperature: ☐ Air ☐ Filter		Checked by:	15
Autorial Composition Land Land Land Land			
CUSTODY SEALS INTACT:			
Cooler □ □ No (Not Intact) □ Not Present	□ N/A		
Sample		Checked by:	<u>896</u>
SAMPLE CONDITION:	Yes	No	N/A
Chain-Of-Custody (COC) document(s) received with samples			
COC document(s) received complete			
☐ Collection date/time, matrix, and/or # of containers logged in based on sample labels			
☐ No analysis requested. ☐ Not relinquished. ☐ No date/time relinquished.			
Sampler's name indicated on COC	. 🗸		
Sample container label(s) consistent with COC			
Sample container(s) intact and good condition			
Proper containers and sufficient volume for analyses requested			
Analyses received within holding time	9		
Aqueous samples received within 15-minute holding time			
☐ pH ☐ Residual Chlorine ☐ Dissolved Sulfides ☐ Dissolved Oxygen	🗆		
Proper preservation noted on COC or sample container	. 🗾		
Unpreserved vials received for Volatiles analysis	>		
Volatile analysis container(s) free of headspace			
Tedlar bag(s) free of condensation CONTAINER TYPE:	🗆		
Solid: 🗆4ozCGJ 🗆8ozCGJ 🗀16ozCGJ 🗀Sleeve () 🗀EnCore	es [®] □Terra	Cores® □	
Aqueous: ♥VOA ♥VOAh □VOAna2 □125AGB □125AGBh □125AGBr	ZÍAGB	□1AGB na ₂ □	1AGB s
□500AGB ☑500AGJ □500AGJs □250AGB □250CGB □250CGB	n 24 1		00PB
□250PB □250PBnu□125PB □125PBznna □100PJ □100PJna ₂ ☑ 2	50Bn 0_		
Air: Tedlar [®] Canister Other: Trip Blank Lot#: 1404020 Container: C: Clear A: Amber P: Plastic G: Glass J: Jar B: Bottle Z: Ziploc/Resealable Bag E: E	Labeled		8m
Preservative: h: HCL n: HNO ₃ na ₂ :Na ₂ S ₂ O ₃ na: NaOH p: H ₃ PO ₄ s: H ₂ SO ₄ u: Ultra-pure znna: ZnAc ₂ +N	•	Scanned by:	A (



Calscience

WORK ORDER #: 14-10- 2 1 3

SAMPLE RECEIPT FORM

Cooler <u>2</u> of <u>2</u>

CLIENT: EST	DATE: _	10/28/	14					
TEMPERATURE: Thermometer ID: SC2 (Criteria: 0.0 °C – 6.0 °C, not froze	n except se	diment/tissue)					
Temperature _ 2 6 _ °C - 0.2 °C (CF) = _ 2 4 _ °C	[™] Blank	☐ Sample						
☐ Sample(s) outside temperature criteria (PM/APM contacted by:)								
☐ Sample(s) outside temperature criteria but received on ice/chilled on same day of sampling.								
☐ Received at ambient temperature, placed on ice for transport by Co								
Ambient Temperature: ☐ Air ☐ Filter		Checked by	: 15					
CUSTODY SEALS INTACT:			_					
Cooler □ □ No (Not Intact) □ Not Present	□ N/A	Checked by:	: <u>LS</u>					
□ Sample □ □ No (Not Intact) □ Not Present		Checked by:	836					
SAMPLE CONDITION:	Yes	No	N/A					
Chain-Of-Custody (COC) document(s) received with samples								
COC document(s) received complete	_							
☐ Collection date/time, matrix, and/or # of containers logged in based on sample labels								
☐ No analysis requested. ☐ Not relinquished. ☐ No date/time relinquished.								
Sampler's name indicated on COC	. 🗷							
Sample container label(s) consistent with COC	. 🗹							
Sample container(s) intact and good condition	. 🗹							
Proper containers and sufficient volume for analyses requested	. 🗹							
Analyses received within holding time	. 🗹							
Aqueous samples received within 15-minute holding time								
☐ pH ☐ Residual Chlorine ☐ Dissolved Sulfides ☐ Dissolved Oxygen	🗆							
Proper preservation noted on COC or sample container	. 🖊							
☐ Unpreserved vials received for Volatiles analysis								
Volatile analysis container(s) free of headspace	🗆							
Tedlar bag(s) free of condensation CONTAINER TYPE:	🗆							
Solid: □4ozCGJ □8ozCGJ □16ozCGJ □Sleeve () □EnCore	es [®] □Terra	Cores [®] □						
Aqueous: □VOA □VOAh □VOAna₂ □125AGB □125AGBh □125AGBr	o Ø1AGB [□1AGB na ₂ □	1AGB s					
□500AGB ☑500AGJ □500AGJs □250AGB □250CGB □250CGB	s □1PB	□1PB na □5	500PB					
□250PB №250PBn □125PB □125PB znna □100PJ □100PJ na ₂ □_								
Air: ☐Tedlar® ☐Canister Other: ☐ Trip Blank Lot#:		/Checked by: . Reviewed by: _	Sm.					
Container: C: Clear A: Amber P: Plastic G: Glass J: Jar B: Bottle Z: Ziploc/Resealable Bag E: E Preservative: h: HCL n: HNO ₃ na ₂ :Na ₂ S ₂ O ₃ na: NaOH p: H ₃ PO ₄ s: H ₂ SO ₄ u: Ultra-pure znna: ZnAc ₂ +N ₃	*	-	8nt					

RAW DATA SHEET FOR METHOD: EPA 8015B (M)

WORK ORDER: INSTRUMENT:

EXTRACTION:

14-10-2131

GC 45

EPA 3510C

D/T EXTRACTED:

2014-10-30 00:00

682 **ANALYZED BY:**

D/T ANALYZED: 2014-10-31 09:41

628 **REVIEWED BY:**

D/T REVIEWED:

2014-11-04 17:13

DATA FILE:

W:\GC 45\GC 45 DATA\2014\141030\14103063.D\14103063

CLIENT SAMPLE NUMBER: ES113

LCS/MB BATCH:

141030B06

SAMPLE VOLUME / WEIGHT:

DEFAULT: 500.00 ml / ACTUAL: 500.00 ml

DEFAULT: 5.00 ml / ACTUAL: 2.50 ml

MS/MSD BATCH: **UNITS:**

ug/L

FINAL VOLUME / WEIGHT: ADJUSTMENT RATIO TO PF:

0.50

COMMENT:

Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are

qualified with a "J" flag. TPH as Diesel is quantified in the carbon range C10-C28.

COMPOUND TPH as Diesel INI. CONC

23400

DF 1.00

CONC 117

DL 11

LOD

12

QUAL

b

LOQ

25

Area Percent Report

Data File Name : W:\GC_45\GC 45 DATA\2014\141030\14103063.D

Page Number : 1

Operator : 682 Vial Number : Vial 63

Instrument : GC 45 Injection Number : 1 Sample Name : 14-10-2131-2 Sequence Line : 63

Instrument Method: C:\CHEM32\1\METHODS\ ->

Acquired on : 31 Oct 14 09:41 am

Report Created on: 04 Nov 14 02:48 pm Analysis Method : 8015B.MTH

Sig. 1 in $W:\GC_45\GC 45\DATA\2014\141030\->$

Pk	Ret Time	Area	Height	Peak	Width	Response %
1	2.024	2.51		VV	0.035	0.239
2	2.092	1.31	0	VV	0.048	0.125
3	2.223	7.06	2	VV	0.039	0.671
4	2.285	1.69	1	VV	0.022	0.161
5	2.334	3.26	1	VV	0.034	0.309
6	2.463	10.00	3	VV	0.048	0.950
7	2.511	3.68	2	VV	0.028	0.349
8	2.638	13.63	3	VV	0.070	1.294
9	2.753	24.19	5	VV	0.060	2.297
10	2.883	24.76	7	VV	0.045	2.351
11	2.943	15.15	5	VV	0.044	1.439
12	3.012	16.40	4	VV	0.062	1.558
13	3.094	8.03	4	VV	0.032	0.762
14	3.168	19.46	4	VV	0.056	1.848
15	3.306	37.68	4	VV	0.109	3.578
16	3.438	14.10	4	VV	0.050	1.339
17	3.508	16.84	4	VV	0.059	1.599
18	3.559	10.28	4	VV	0.041	0.976
19	3.602	13.03	3	VV	0.048	1.238
20	3.715	25.80	4	VV	0.089	2.450
21	3.849	7.80	3	VV	0.040	0.741
22	3.911	9.29	2	VV	0.067	0.882
23	4.017	11.82	3	VV	0.052	1.123
24	4.171	1.93	1	VV	0.034	0.183
25	4.238	5.50	1	VV	0.075	0.523
26	4.399	7.57	1	VV	0.119	0.718
27	4.593	1.54	0	VV	0.046	0.146
28	4.690	5.13	1	VV	0.090	0.487
29	4.848	2.51	1	VV	0.046	0.239
30	4.909	1.46	1	VV	0.036	0.138
31	4.967	1.29	1	VV	0.034	0.122
32	5.069	3.16	1	VV	0.044	0.300
33	5.220	2.72	1	VV	0.044	0.258
34	5.278	1.24	0	VV	0.044	0.118
35	5.373	721.21	503	VV	0.024	68.488

Area Percent Report

Data File Name : W:\GC_45\GC 45 DATA\2014\141030\14103063.D

Page Number : 2

Operator : 682 Vial Number : Vial 63

Instrument : GC 45 Injection Number : 1 Sample Name : 14-10-2131-2 Sequence Line : 63

Instrument Method: C:\CHEM32\1\METHODS\ ->

Acquired on : 31 Oct 14 09:41 am

Report Created on: 04 Nov 14 02:48 pm Analysis Method : 8015B.MTH

Software Revision: Rev. B.03.02 [341] Copyright © Agilent Technologies

1053.05

Area Percent Report

Data File Name : W:\GC 45\GC 45 DATA\2014\141030\14103063.D

Page Number : 3

Operator : 682

Vial Number : Vial 63

Instrument

: GC 45

Injection Number: 1

Sample Name

: 14-10-2131-2

Sequence Line : 63

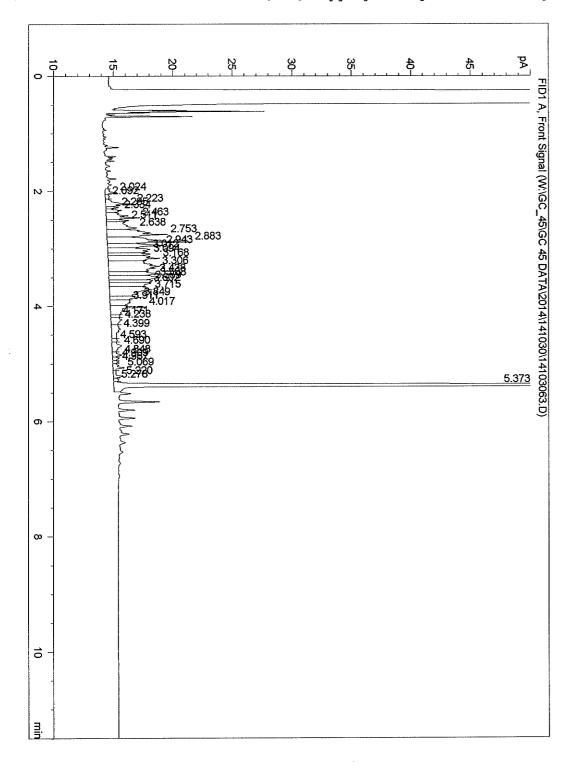
Instrument Method: C:\CHEM32\1\METHODS\

Acquired on

: 31 Oct 14

09:41 am

Report Created on: 04 Nov 14 02:48 pm Analysis Method: 8015B.MTH



RAW DATA SHEET FOR METHOD: EPA 8015B (M)

WORK ORDER:

14-10-2131

INSTRUMENT:

GC 45

EXTRACTION:

EPA 3510C

D/T EXTRACTED: 2014-10-30 00:00

ANALYZED BY:

682 D/T ANALYZED:

2014-10-31 09:59

628 **REVIEWED BY:**

D/T REVIEWED:

2014-11-04 17:13

DATA FILE:

W:\GC 45\GC 45 DATA\2014\141030\14103064.D\14103064

CLIENT SAMPLE NUMBER: ES114

LCS/MB BATCH: MS/MSD BATCH:

141030B06 141030S06 SAMPLE VOLUME / WEIGHT:

DEFAULT: 500.00 ml / ACTUAL: 500.00 ml DEFAULT: 5.00 ml / ACTUAL: 2.50 ml

FINAL VOLUME / WEIGHT:

UNITS:

ug/L

ADJUSTMENT RATIO TO PF:

0.50

COMMENT:

Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are

qualified with a "J" flag. TPH as Diesel is quantified in the carbon range C10-C28.

COMPOUND

INI. CONC DF 409000

CONC

<u>DL</u>

LOD

QUAL

TPH as Diesel

1.00

2050

11

12

b

LOQ

25

Area Percent Report

-

Data File Name : W:\GC_45\GC 45 DATA\2014\141030\14103064.D

Page Number : 1

Operator : 682 Vial Number : Vial 64

Instrument : GC 45 Injection Number : 1
Sample Name : 14-10-2131-3 Sequence Line : 64

Instrument Method: C:\CHEM32\1\METHODS\ ->

Acquired on : 31 Oct 14 09:59 am

Report Created on: 04 Nov 14 02:49 pm Analysis Method : 8015B.MTH

Sig. 1 in W:\GC 45\GC 45 DATA\2014\141030\ ->

Pk	Ret Time	Area	Height			Response %
				1		
1	1.947	36.87	21	VV	0.025	0.554
2	2.028	166.85		VV	0.032	2.505
3	2.173	220.62	128		0.027	3.313
4	2.258	188.02	81	VV	0.032	2.823
5	2.353	227.05	110	VV	0.029	3.409
6	2.457	789.99	535		0.022	11.862
7	2.544	191.88		VV	0.053	2.881
8	2.600	106.50	52	VV	0.030	1.599
9	2.642	116.77		VV	0.032	1.753
10	2.729	439.52	208		0.030	6.599
11	2.769	366.68	248		0.022	5.506
12	2.853	424.12	97	VV	0.056	6.368
13	2.948	289.07		VV	0.050	4.340
14	3.014	234.69	134	VV	0.025	3.524
15	3.058	266.23	81	VV	0.043	3.998
16	3.146	250.38	51	VV	0.064	3.759
17	3.298	368.20	51	VV	0.087	5.529
18	3.431	256.38		VV	0.083	3.850
19	3.498	130.37	37	VV	0.045	1.957
20	3.558	234.19		VV	0.074	3.516
21	3.687	134.10		VV	0.053	2.013
22	3.773	105.43		VV	0.060	1.583
23	3.853	46.10		VV	0.035	0.692
24	3.899	37.48		VV	0.032	0.563
25	3.938	58.23		VV	0.042	0.874
26	4.010	116.07		VV	0.102	1.743
27	4.261	19.12		VV	0.056	0.287
28	4.348	45.09		VV	0.154	0.677
29	4.681	20.80		VV	0.098	0.312
30	4.832	7.60		VV	0.050	0.114
31	4.891	5.24		VV	0.043	0.079
32	4.950	5.08		VV	0.046	0.076
33	5.048	12.74		VV	0.082	0.191
34	5.199	6.89		VV	0.054	0.103
35	5.250	4.12		VV	0.043	0.062
36	5.351	731.57	493	VV	0.024	10.984

Area Percent Report

Data File Name : W:\GC_45\GC 45 DATA\2014\141030\14103064.D

Page Number : 2

Vial Number : Vial 64 : 682 Operator

: GC 45 Instrument

Injection Number: 1 Sequence Line : 64 Sample Name : 14-10-2131-3

Instrument Method: C:\CHEM32\1\METHODS\ ->

Acquired on : 31 Oct 14 09:59 am

Analysis Method : 8015B.MTH Report Created on: 04 Nov 14 02:49 pm

Software Revision: Rev. B.03.02 [341] Copyright © Agilent Technologies

Total area = 6660.03

Area Percent Report

Data File Name : W:\GC 45\GC 45 DATA\2014\141030\14103064.D

Page Number : 3

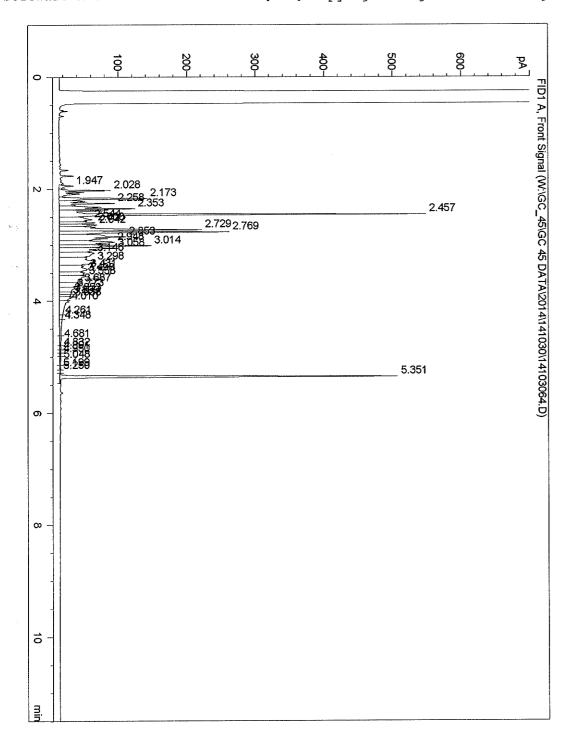
Operator : 682 Vial Number : Vial 64

Instrument : GC 45 Injection Number : 1 Sample Name : 14-10-2131-3 Sequence Line : 64

Instrument Method: C:\CHEM32\1\METHODS\ -:

Acquired on : 31 Oct 14 09:59 am

Report Created on: 04 Nov 14 02:49 pm Analysis Method : 8015B.MTH



RAW DATA SHEET FOR METHOD: EPA 8015B (M)

WORK ORDER: 14-10-2131 ANALYZED BY: 682

INSTRUMENT: GC 45 D/T ANALYZED: 2014-10-31 10:41

EXTRACTION: EPA 3510C REVIEWED BY: 628

<u>D/T EXTRACTED:</u> 2014-10-30 00:00 <u>D/T REVIEWED:</u> 2014-11-04 17:13

DATA FILE: W:\GC 45\GC 45 DATA\2014\141030\14103065.D\14103065

4 <u>CLIENT SAMPLE NUMBER:</u> ES115

LCS/MB BATCH: 141030B06 SAMPLE VOLUME / WEIGHT: DEFAULT: 500.00 ml / ACTUAL: 500.00 ml

MS/MSD BATCH: DEFAULT: 5.00 ml / ACTUAL: 2.50 ml

UNITS: ug/L ADJUSTMENT RATIO TO PF: 0.50

COMMENT: Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are

qualified with a "J" flag. TPH as Diesel is quantified in the carbon range C10-C28.

LOQ QUAL LOD CONC DL INI. CONC DF COMPOUND 25 b 2030 11 12 406000 1.00 TPH as Diesel

Area Percent Report

Data File Name : W:\GC_45\GC 45 DATA\2014\141030\14103065.D

Page Number : 1

Operator : 682 Vial Number : Vial 65

Instrument : GC 45 Injection Number : 1 Sample Name : 14-10-2131-4 Sequence Line : 65

Instrument Method: C:\CHEM32\1\METHODS\ ->

Acquired on : 31 Oct 14 10:41 am

Report Created on: 04 Nov 14 02:50 pm Analysis Method : 8015B.MTH

Sig. 1 in W:\GC_45\GC 45 DATA\2014\141030\ ->

Pk	Ret Time	Area	Height	Peak	Width	Response %
					•	-
1	1.946	36.71	21	VV	0.026	0.555
2	2.028	169.15	75	VV	0.031	2.559
3	2.173	226.98	127	VV	0.028	3.434
4	2.222	61.36	36	VV	0.025	0.928
5	2.257	128.76	81	VV	0.024	1.948
6	2.352	231.89	109	VV	0.030	3.508
7	2.457	829.85	551		0.023	12.555
8	2.542	189.89	47	VV	0.055	2.873
9	2.600	104.99	51	VV	0.030	1.588
10	2.641	115.18	56	VV	0.030	1.743
11	2.730	449.33	212		0.030	6.798
12	2.770	375.36	260	VV	0.022	5.679
13	2.852	404.48	91	VV	0.057	6.119
14	2.947	210.06	78	VV	0.041	3.178
15	3.014	316.74	136	VV	0.032	4.792
16	3.059	267.38	81	VV	0.044	4.045
17	3.149	231.23	48	VV	0.063	3.498
18	3.296	424.16	48	VV	0.113	6.417
19	3.431	166.71	41	VV	0.053	2.522
20	3.502	122.86	33	VV	0.047	1.859
21	3.561	218.09	37	VV	0.075	3.299
22	3.690	126.19	30	VV	0.054	1.909
23	3.783	94.82	21	VV	0.059	1.434
24	3.861	40.17		VV	0.033	
25	3.907	37.37		VV	0.039	
26	3.950	52.52		VV	0.043	
27	4.023	152.63		VV	0.159	
28	4.603	7.19		VV	0.051	
29	4.719	16.45		VV	0.095	
30	4.871	7.28		VV	0.051	
31	4.931	5.14		VV	0.054	
32	4.994	4.19		VV	0.043	
33	5.096	12.49		VV	0.083	
34	5.252	5.63		VV	0.045	
35	5.313	3.60		VV	0.044	
36	5.408	762.98	520	VV	0.023	11.543

Area Percent Report

Data File Name : W:\GC_45\GC 45 DATA\2014\141030\14103065.D

Page Number : 2

Operator : 682 Vial Number : Vial 65

Instrument : GC 45 Injection Number : 1 Sample Name : 14-10-2131-4 Sequence Line : 65

Instrument Method: C:\CHEM32\1\METHODS\ ->

Acquired on : 31 Oct 14 10:41 am

Report Created on: 04 Nov 14 02:50 pm Analysis Method : 8015B.MTH

Software Revision: Rev. B.03.02 [341] Copyright © Agilent Technologies

Total area = 6609.83

Area Percent Report

Data File Name : W:\GC 45\GC 45 DATA\2014\141030\14103065.D

Page Number : 3

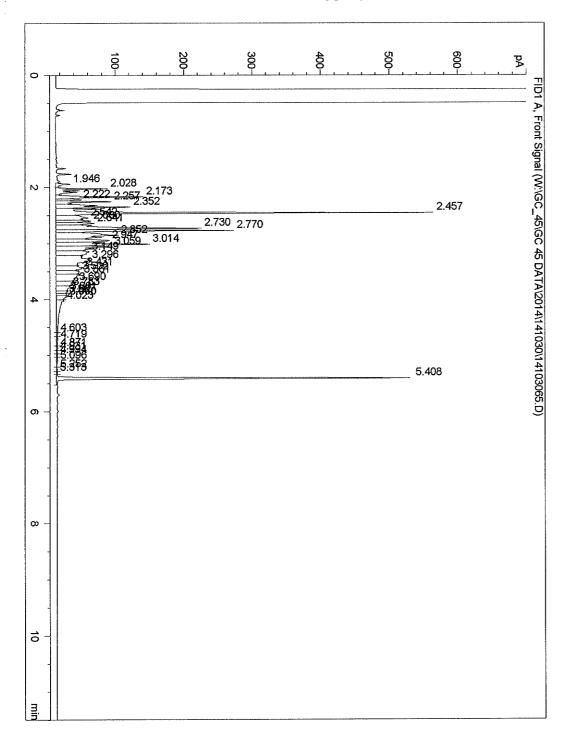
Operator : 682 Vial Number : Vial 65

Instrument : GC 45 Injection Number : 1
Sample Name : 14-10-2131-4 Sequence Line : 65

Instrument Method: C:\CHEM32\1\METHODS\ -

Acquired on : 31 Oct 14 10:41 am

Report Created on: 04 Nov 14 02:50 pm Analysis Method : 8015B.MTH



RAW DATA SHEET FOR METHOD: EPA 8015B (M)

WORK ORDER:

14-10-2131

INSTRUMENT: GC 45

EPA 3510C

EXTRACTION: D/T EXTRACTED: 2014-10-30 00:00

ANALYZED BY:

682

2014-10-31 10:58 D/T ANALYZED:

REVIEWED BY: 628

LOD

12

2014-11-04 17:13 D/T REVIEWED:

DATA FILE:

W:\GC 45\GC 45 DATA\2014\141030\14103066.D\14103066

CLIENT SAMPLE NUMBER: ES116

LCS/MB BATCH:

141030B06 141030S06 SAMPLE VOLUME / WEIGHT:

DEFAULT: 500.00 ml / ACTUAL: 500.00 ml

DEFAULT: 5.00 ml / ACTUAL: 2.50 ml FINAL VOLUME / WEIGHT:

MS/MSD BATCH: **UNITS:**

ug/L

ADJUSTMENT RATIO TO PF:

0.50

COMMENT:

Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are

qualified with a "J" flag. TPH as Diesel is quantified in the carbon range C10-C28.

COMPOUND TPH as Diesel INI. CONC DF 16000

CONC 79.9 1.00

DL 11

LOQ 25

QUAL b

Area Percent Report

Data File Name : W:\GC_45\GC 45 DATA\2014\141030\14103066.D

Page Number : 1

Operator : 682 Vial Number : Vial 66

Instrument : GC 45 Injection Number : 1
Sample Name : 14-10-2131-5 Sequence Line : 66

Instrument Method: C:\CHEM32\1\METHODS\ ->

Acquired on : 31 Oct 14 10:58 am

Report Created on: 04 Nov 14 02:51 pm Analysis Method : 8015B.MTH

Software Revision: Rev. B.03.02 [341] Copyright © Agilent Technologies

Sig. 1 in W:\GC 45\GC 45 DATA\2014\141030\ \rightarrow

Pk	Ret Time	Area	Height		Width	Response 🖇
		•		•	•	
1	2.219	0.45	0	VV	0.038	0.060
2	2.270	0.27	0	VV	0.031	0.036
3	2.457	0.84	0	VV	0.065	0.113
4	2.634	1.87	1	VV	0.049	0.252
5	2.750	4.97	1	VV	0.064	0.669
6	2.837	1.49	1	VV	0.037	0.200
7	2.884	1.51	1	VV	0.032	0.204
8	2.950	2.03	1	VV	0.045	0.273
9	3.169	12.58	2	VV	0.078	1.695
10	3.306	10.78	2	VV	0.070	1.452
11	3.372	8.28	2	VV	0.058	1.115
12	3.436	8.46	3	VV	0.046	1.139
13	3.511	26.81	3	VV	0.117	3.611
14	3.701	20.15	2	VV	0.122	2.714
15	3.846	7.00	2	VV	0.051	0.943
16	3.937	5.92	2	VV	0.052	0.797
17	4.012	7.26	3	VV	0.036	0.978
18	4.111	8.62	2	VV	0.075	1.161
19	4.228	13.65	2	VV	0.096	1.839
20	4.387	20.35	2	VV	0.137	2.740
21	4.560	13.57	2	VV	0.100	1.828
22	4.662	12.72	2	VV	0.090	1.714
23	4.826	7.22	2	VV	0.064	0.972
24	4.889	3.85	1	VV	0.042	0.518
25	4.946	6.18	1	VV	0.061	0.832
26	5.043	8.15	1	VV	0.074	1.097
27	5.193	3.65	1	VV	0.041	0.492
28	5.249	2.50	1	VV	0.048	0.337
29	5.344	521.35	374	VV	0.022	70.220

Total area = 742.45

Area Percent Report

Data File Name : W:\GC_45\GC 45 DATA\2014\141030\14103066.D

Page Number

: 2

: 682

Operator Instrument

: GC 45

Sample Name

: 14-10-2131-5

Vial Number

: Vial 66

Injection Number: 1

Sequence Line

Instrument Method: C:\CHEM32\1\METHODS\

: 66

Acquired on

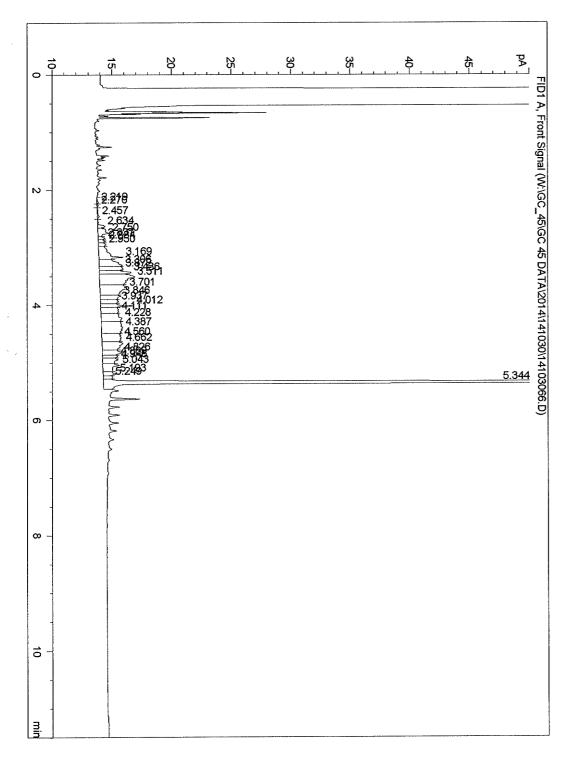
: 31 Oct 14

10:58 am

Report Created on: 04 Nov 14

02:51 pm

Analysis Method : 8015B.MTH



Area Percent Report

Data File Name : W:\GC_45\GC 45 DATA\2014\141030\14103051.D

Page Number : 3

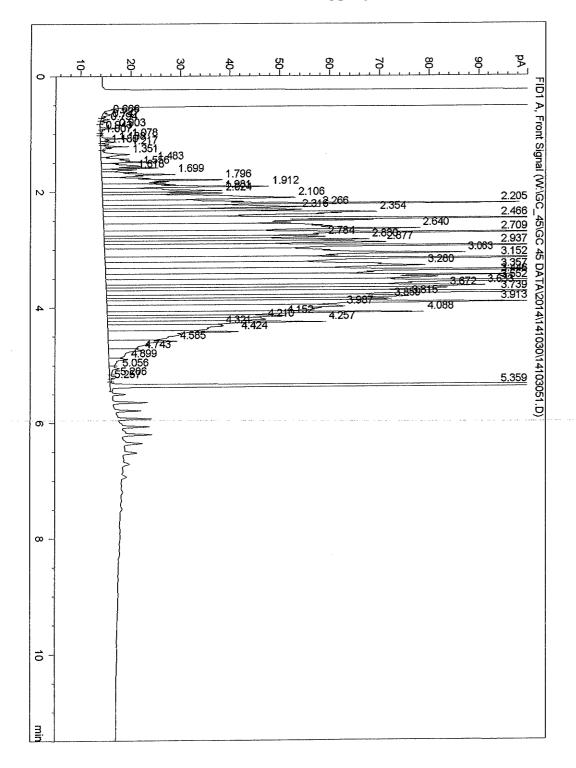
Operator : 682 Vial Number : Vial 51

Instrument : GC 45 Injection Number : 1
Sample Name : D400 C28 50 L080814D Sequence Line : 51

Instrument Method: C:\CHEM32\1\METHODS\ ->

Acquired on : 31 Oct 14 06:11 am

Report Created on: 04 Nov 14 01:18 pm Analysis Method : 8015B.MTH





Calscience



WORK ORDER NUMBER: 14-10-2234

The difference is service



AIR | SOIL | WATER | MARINE CHEMISTRY

Analytical Report For

Client: Environmental Science International, Inc.

Client Project Name: Red Hill LTM 112066

Attention: Jeff Hattemer

354 Uluniu Street, Suite 304 Kailua, HI 96734-2500

Richard Vellas

Approved for release on 11/05/2014 by:

Richard Villafania Project Manager



ResultLink >

Email your PM >

Eurofins Calscience, Inc. (Calscience) certifies that the test results provided in this report meet all NELAC requirements for parameters for which accreditation is required or available. Any exceptions to NELAC requirements are noted in the case narrative. The original report of subcontracted analyses, if any, is attached to this report. The results in this report are limited to the sample(s) tested and any reproduction thereof must be made in its entirety. The client or recipient of this report is specifically prohibited from making material changes to said report and, to the extent that such changes are made, Calscience is not responsible, legally or otherwise. The client or recipient agrees to indemnify Calscience for any defense to any litigation which may arise.



Contents

Client Project Name: Red Hill LTM 112066 Work Order Number: 14-10-2234

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5	Glossary of Terms and Qualifiers	28
6	Chain-of-Custody/Sample Receipt Form	29



Work Order Narrative

Work Order: 14-10-2234 Page 1 of 1

Condition Upon Receipt:

Samples were received under Chain-of-Custody (COC) on 10/29/14. They were assigned to Work Order 14-10-2234.

Unless otherwise noted on the Sample Receiving forms all samples were received in good condition and within the recommended EPA temperature criteria for the methods noted on the COC. The COC and Sample Receiving Documents are integral elements of the analytical report and are presented at the back of the report.

Holding Times:

All samples were analyzed within prescribed holding times (HT) and/or in accordance with the Calscience Sample Acceptance Policy unless otherwise noted in the analytical report and/or comprehensive case narrative, if required.

Any parameter identified in 40CFR Part 136.3 Table II that is designated as "analyze immediately" with a holding time of <= 15 minutes (40CFR-136.3 Table II, footnote 4), is considered a "field" test and the reported results will be qualified as being received outside of the stated holding time unless received at the laboratory within 15 minutes of the collection time.

Quality Control:

All quality control parameters (QC) were within established control limits except where noted in the QC summary forms or described further within this report.

Additional Comments:

Air - Sorbent-extracted air methods (EPA TO-4A, EPA TO-10, EPA TO-13A, EPA TO-17): Analytical results are converted from mass/sample basis to mass/volume basis using client-supplied air volumes.

New York NELAP air certification does not certify for all reported methods and analytes, reference the accredited items here: http://www.calscience.com/PDF/New_York.pdf

Solid - Unless otherwise indicated, solid sample data is reported on a wet weight basis, not corrected for % moisture. All QC results are always reported on a wet weight basis.

Subcontractor Information:

Unless otherwise noted below (or on the subcontract form), no samples were subcontracted.



Environmental Science International, Inc. Date Received: 10/29/14 Work Order: 14-10-2234 354 Uluniu Street, Suite 304 Kailua, HI 96734-2500 Preparation: **EPA 3510C** Method: EPA 8015B (M) Units: ug/L

Project: Red Hill LTM 112066 Page 1 of 1

Client Sample Number	Lab Sample Number	Date/Time Collected	Matrix	Instrument	Date Prepared	Date/Time Analyzed	QC Batch ID
ES117	14-10-2234-2-G	10/28/14 12:00	Aqueous	GC 45	10/30/14	10/31/14 11:16	141030B06

Comment(s): - Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are qualified with a "J" flag.

- TPH as Diesel is quantified in the carbon range C10-C28.

<u>DF</u> **Parameter** Result <u>DL</u> <u>LOD</u> <u>LOQ</u> **Qualifiers** TPH as Diesel HD,J 11 12 25 1.00 22

Control Limits Qualifiers <u>Surrogate</u> Rec. (%)

n-Octacosane 85 51-141

E	ES118	14-10-2234-3-G	10/28/14 14:30	Aqueous	GC 45	10/30/14	10/31/14 11:35	141030B06
			17.00				11.00	

- Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are qualified with a "J" flag. Comment(s):

- TPH as Diesel is quantified in the carbon range C10-C28.

<u>DF</u> <u>Parameter</u> Result LOD LOQ Qualifiers TPH as Diesel 16 11 12 25 1.00 HD,J

Surrogate Rec. (%) **Control Limits** Qualifiers

51-141 n-Octacosane 70

Method Blank	C	099-15-516-206	N/A Aq	ueous	GC 45	10/30/14	10/31/14 07:03	141030B06
Comment(s):	- Results were evaluated to the	ne MDL (DL), conce	entrations >= to the I	MDL (DL) but < RL	(LOQ), if found, a	are qualified with a	a "J" flag.
<u>Parameter</u>		<u>Result</u>	<u>DL</u>	<u>LOD</u>		<u>LOQ</u>	<u>DF</u>	<u>Qualifiers</u>
TPH as Diesel		<12	11	12		25	1.00	U
Surrogate n-Octacosane		<u>Rec. (%)</u> 117	Control Limits 51-141	Qualifi	<u>ers</u>			



10/29/14

14-10-2234



Kailua, HI 96734-2500

Analytical Report

Environmental Science International, Inc. Date Received: 354 Uluniu Street, Suite 304 Work Order:

Preparation:

EPA 3005A Filt. Method:

EPA 6020 Units: ug/L

Page 1 of 1 Project: Red Hill LTM 112066

Client Sample N	Number	Lab Sample Number	Date/Time Collected	Matrix	Instrument	Date Prepared	Date/Time Analyzed	QC Batch ID
ES118		14-10-2234-3-J	10/28/14 14:30	Aqueous	ICP/MS 04	10/31/14	11/03/14 15:10	141031L04F
Comment(s):	- Results were evaluated to	the MDL (DL), cond	centrations >= to	the MDL (DL	_) but < RL (LO	Q), if found, a	re qualified with a	a "J" flag.
<u>Parameter</u>		Result	<u>DL</u>	<u>LOD</u>	<u>LO</u>	Q	<u>DF</u>	<u>Qualifiers</u>
Lead		<0.200	0.0898	0.200	1.0	0	1.00	U

Method Blank	099-14-497-10	4 N/A	Aqueous	ICP/MS 04	10/31/14 11/0 14:4	03/14 141031L04F 48
Comment(s):	- Results were evaluated to the MDL (DL),	concentrations	>= to the MDL (DL)	but < RL (LOQ)	, if found, are qualifi	ed with a "J" flag.
<u>Parameter</u>	Result	<u>DL</u>	<u>LOD</u>	<u>LOQ</u>	<u>DF</u>	<u>Qualifiers</u>
Lead	<0.200	0.089	98 0.200	1.00	1.00	U





Environmental Science International, Inc. 354 Uluniu Street, Suite 304

Kailua, HI 96734-2500

Date Received:

10/29/14 14-10-2234

Work Order: Preparation:

EPA 3510C

Method:

EPA 8270C SIM PAHs

Units:

ug/L

Project: Red Hill LTM 112066

Page 1 of 3

Client Sample Number	Lab Sample Number	Date/Time Collected	Matrix	Instrument	Date Prepared	Date/Time Analyzed	QC Batch ID
ES117	14-10-2234-2-I	10/28/14 12:00	Aqueous	GC/MS AAA	10/30/14	11/03/14 17:03	141030L08
Comment(s): - Results were evaluated	to the MDL (DL), con	centrations >= to tl	he MDL (DL) but < RL (LOC	(), if found, are	qualified with a	a "J" flag.
<u>Parameter</u>	<u>Result</u>	<u>DL</u>	<u>LOD</u>	LOC	<u>]</u>	<u>DF</u>	<u>Qualifiers</u>
Naphthalene	< 0.049	0.033	0.049	0.19) 1	.00	U
2-Methylnaphthalene	< 0.049	0.045	0.049	0.19) 1	.00	U
1-Methylnaphthalene	<0.097	0.050	0.097	0.19) 1	.00	U
Acenaphthylene	<0.049	0.043	0.049	0.19) 1	.00	U
Acenaphthene	< 0.049	0.026	0.049	0.19) 1	.00	U
Fluorene	< 0.049	0.041	0.049	0.19) 1	.00	U
Phenanthrene	< 0.049	0.026	0.049	0.19) 1	.00	U
Anthracene	< 0.049	0.028	0.049	0.19) 1	.00	U
Fluoranthene	< 0.049	0.045	0.049	0.19) 1	.00	U
Pyrene	< 0.049	0.020	0.049	0.19) 1	.00	U
Benzo (a) Anthracene	<0.049	0.032	0.049	0.19) 1	.00	U
Chrysene	< 0.049	0.024	0.049	0.19) 1	.00	U
Benzo (k) Fluoranthene	<0.049	0.030	0.049	0.19) 1	.00	U
Benzo (b) Fluoranthene	<0.049	0.017	0.049	0.19) 1	.00	U
Benzo (a) Pyrene	<0.049	0.022	0.049	0.19) 1	.00	U
Indeno (1,2,3-c,d) Pyrene	<0.049	0.021	0.049	0.19) 1	.00	U
Dibenz (a,h) Anthracene	< 0.049	0.046	0.049	0.19) 1	.00	U
Benzo (g,h,i) Perylene	<0.097	0.080	0.097	0.19) 1	.00	U
Surrogate	Rec. (%)	Control Limi	its Qualif	<u>iers</u>			
Nitrobenzene-d5	61	28-139					
2-Fluorobiphenyl	81	33-144					
p-Terphenyl-d14	83	23-160					



Environmental Science International, Inc. 354 Uluniu Street, Suite 304

Kailua, HI 96734-2500

Date Received:

10/29/14 14-10-2234

Work Order: Preparation:

EPA 3510C

Method:

EPA 8270C SIM PAHs

Units:

ug/L

Project: Red Hill LTM 112066

Page 2 of 3

Client Sample Number	Lab Sample Number	Date/Time Collected	Matrix	Instrument	Date Prepared	Date/Time Analyzed	QC Batch ID
ES118	14-10-2234-3-I	10/28/14 14:30	Aqueous	GC/MS AAA	10/30/14	11/03/14 17:26	141030L08
Comment(s): - Results were evaluated	to the MDL (DL), con	centrations >= to t	he MDL (DL	.) but < RL (LOC	Q), if found, a	are qualified with a	a "J" flag.
<u>Parameter</u>	<u>Result</u>	<u>DL</u>	LOD	LOC	<u>2</u>	<u>DF</u>	<u>Qualifiers</u>
Naphthalene	<0.048	0.033	0.048	0.19	9	1.00	U
2-Methylnaphthalene	<0.048	0.045	0.048	0.19	9	1.00	U
1-Methylnaphthalene	<0.096	0.050	0.096	0.19	9	1.00	U
Acenaphthylene	<0.048	0.043	0.048	0.19	9	1.00	U
Acenaphthene	<0.048	0.026	0.048	0.19	9	1.00	U
Fluorene	<0.048	0.041	0.048	0.19	9	1.00	U
Phenanthrene	<0.048	0.026	0.048	0.19	9	1.00	U
Anthracene	<0.048	0.028	0.048	0.19	9	1.00	U
Fluoranthene	<0.048	0.045	0.048	0.19	9	1.00	U
Pyrene	<0.048	0.020	0.048	0.19	9	1.00	U
Benzo (a) Anthracene	<0.048	0.031	0.048	0.19	9	1.00	U
Chrysene	<0.048	0.024	0.048	0.19	9	1.00	U
Benzo (k) Fluoranthene	<0.048	0.030	0.048	0.19	9	1.00	U
Benzo (b) Fluoranthene	<0.048	0.017	0.048	0.19	9	1.00	U
Benzo (a) Pyrene	<0.048	0.021	0.048	0.19	9	1.00	U
Indeno (1,2,3-c,d) Pyrene	<0.048	0.020	0.048	0.19	9	1.00	U
Dibenz (a,h) Anthracene	<0.048	0.046	0.048	0.19	9	1.00	U
Benzo (g,h,i) Perylene	<0.096	0.079	0.096	0.19	9	1.00	U
<u>Surrogate</u>	Rec. (%)	Control Limi	<u>its</u> Qualif	<u>iers</u>			
Nitrobenzene-d5	62	28-139					
2-Fluorobiphenyl	88	33-144					
p-Terphenyl-d14	94	23-160					



Environmental Science International, Inc.

Kailua, HI 96734-2500

Date Received: Work Order:

10/29/14 14-10-2234

354 Uluniu Street, Suite 304

Preparation:

EPA 3510C

Method:

EPA 8270C SIM PAHs

Units:

ug/L

Project: Red Hill LTM 112066

Page 3 of 3

Client Sample Number	Lab Sample Number	Date/Time Collected	Matrix	Instrument	Date Prepared	Date/Time Analyzed	QC Batch ID
Method Blank	099-15-148-67	N/A	Aqueous	GC/MS AAA	10/30/14	10/31/14 17:17	141030L08
Comment(s): - Results were evaluated	to the MDL (DL), cor	ncentrations >= to the	he MDL (DL) but < RL (LOC	Q), if found, a	are qualified with	a "J" flag.
<u>Parameter</u>	Result	<u>DL</u>	<u>LOD</u>	LOC	<u>)</u>	<u>DF</u>	<u>Qualifiers</u>
Naphthalene	< 0.050	0.034	0.050	0.20)	1.00	U
2-Methylnaphthalene	< 0.050	0.046	0.050	0.20)	1.00	U
1-Methylnaphthalene	<0.10	0.052	0.10	0.20)	1.00	U
Acenaphthylene	< 0.050	0.045	0.050	0.20)	1.00	U
Acenaphthene	< 0.050	0.027	0.050	0.20)	1.00	U
Fluorene	< 0.050	0.043	0.050	0.20)	1.00	U
Phenanthrene	< 0.050	0.027	0.050	0.20)	1.00	U
Anthracene	< 0.050	0.029	0.050	0.20)	1.00	U
Fluoranthene	< 0.050	0.047	0.050	0.20)	1.00	U
Pyrene	< 0.050	0.020	0.050	0.20)	1.00	U
Benzo (a) Anthracene	< 0.050	0.033	0.050	0.20)	1.00	U
Chrysene	< 0.050	0.025	0.050	0.20)	1.00	U
Benzo (k) Fluoranthene	< 0.050	0.031	0.050	0.20)	1.00	U
Benzo (b) Fluoranthene	< 0.050	0.018	0.050	0.20)	1.00	U
Benzo (a) Pyrene	< 0.050	0.022	0.050	0.20)	1.00	U
Indeno (1,2,3-c,d) Pyrene	< 0.050	0.021	0.050	0.20)	1.00	U
Dibenz (a,h) Anthracene	< 0.050	0.048	0.050	0.20)	1.00	U
Benzo (g,h,i) Perylene	<0.10	0.082	0.10	0.20)	1.00	U
Surrogate	Rec. (%)	Control Limi	its Qualifi	ers			
Nitrobenzene-d5	53	28-139					
2-Fluorobiphenyl	68	33-144					
p-Terphenyl-d14	72	23-160					



Environmental Science International, Inc. 354 Uluniu Street, Suite 304

Kailua, HI 96734-2500

Date Received:

10/29/14 14-10-2234

Work Order: Preparation:

EPA 5030C

Method:

GC/MS / EPA 8260B

Units:

ug/L

Project: Red Hill LTM 112066

Page 1 of 8

Client Sample Number	Lab Sample Number	Date/Time Collected	Matrix	Instrument	Date Date/Ti Prepared Analyze	
ES TRIP	14-10-2234-1-A	10/28/14 11:00	Aqueous	GC/MS OO	10/31/14 11/01/1 01:14	4 141031L047
Comment(s): - Results were evaluated	to the MDL (DL), con	centrations >= 1	to the MDL (DI	_) but < RL (LOC	Q), if found, are qualified	with a "J" flag.
<u>Parameter</u>	<u>Result</u>	<u>DL</u>	<u>LOD</u>	LOC	<u>DF</u>	Qualifiers
Acetone	<10	6.0	10	20	1.00	U,ICH
Benzene	<0.50	0.14	0.50	1.0	1.00	U
Bromodichloromethane	<0.50	0.21	0.50	5.0	1.00	U
Bromoform	<1.0	0.50	1.0	10	1.00	U
Bromomethane	<5.0	3.9	5.0	20	1.00	U,IJ
2-Butanone	<5.0	2.2	5.0	10	1.00	U
Carbon Tetrachloride	<0.50	0.23	0.50	1.0	1.00	U
Chlorobenzene	<0.50	0.17	0.50	5.0	1.00	U
Chloroethane	<5.0	2.3	5.0	10	1.00	U
Chloroform	<0.50	0.46	0.50	5.0	1.00	U
Chloromethane	<2.0	1.8	2.0	10	1.00	U
Dibromochloromethane	<0.50	0.25	0.50	1.0	1.00	U
1,2-Dibromo-3-Chloropropane	<2.0	1.2	2.0	10	1.00	U
1,2-Dibromoethane	<0.50	0.36	0.50	1.0	1.00	U
1,2-Dichlorobenzene	<0.50	0.46	0.50	1.0	1.00	U
1,3-Dichlorobenzene	<0.50	0.40	0.50	1.0	1.00	U
1,4-Dichlorobenzene	0.43	0.43	0.50	1.0	1.00	J
1,1-Dichloroethane	<0.50	0.28	0.50	5.0	1.00	U
1,2-Dichloroethane	<0.50	0.24	0.50	1.0	1.00	U
1,1-Dichloroethene	<0.50	0.43	0.50	1.0	1.00	U
c-1,2-Dichloroethene	<0.50	0.48	0.50	1.0	1.00	U
t-1,2-Dichloroethene	<0.50	0.37	0.50	1.0	1.00	U
1,2-Dichloropropane	<0.50	0.42	0.50	5.0	1.00	U
c-1,3-Dichloropropene	<0.50	0.25	0.50	1.0	1.00	U
t-1,3-Dichloropropene	<0.50	0.25	0.50	1.0	1.00	U
Ethylbenzene	<0.50	0.14	0.50	1.0	1.00	U
Methylene Chloride	<1.0	0.64	1.0	5.0	1.00	U
4-Methyl-2-Pentanone	<5.0	4.4	5.0	10	1.00	U
Styrene	<0.50	0.17	0.50	1.0	1.00	U
1,1,1,2-Tetrachloroethane	<0.50	0.40	0.50	1.0	1.00	U
1,1,2,2-Tetrachloroethane	<0.50	0.41	0.50	1.0	1.00	U
Tetrachloroethene	<0.50	0.39	0.50	5.0	1.00	U
Toluene	<0.50	0.24	0.50	1.0	1.00	U
1,2,4-Trichlorobenzene	<1.0	0.50	1.0	5.0	1.00	U
1,1,1-Trichloroethane	<0.50	0.30	0.50	5.0	1.00	U
Hexachloro-1,3-Butadiene	<0.50	0.32	0.50	1.0	1.00	U



Environmental Science International, Inc. 354 Uluniu Street, Suite 304

Kailua, HI 96734-2500

Date Received: Work Order: Preparation:

Qualifiers

Method: Units:

10/29/14 14-10-2234 EPA 5030C

GC/MS / EPA 8260B

ug/L Page 2 of 8

Project: Red Hill LTM 112066

<u>Parameter</u>	Result	<u>DL</u>	<u>LOD</u>	<u>LOQ</u>	<u>DF</u>	Qualifiers
1,1,2-Trichloroethane	<0.50	0.38	0.50	1.0	1.00	U
Trichloroethene	<0.50	0.37	0.50	1.0	1.00	U
1,2,3-Trichloropropane	<1.0	0.64	1.0	5.0	1.00	U
Vinyl Chloride	<0.50	0.30	0.50	1.0	1.00	U
p/m-Xylene	<1.0	0.30	1.0	10	1.00	U
o-Xylene	<0.50	0.23	0.50	1.0	1.00	U
Methyl-t-Butyl Ether (MTBE)	<0.50	0.31	0.50	1.0	1.00	U
Gasoline Range Organics	<30	26	30	50	1.00	U

Surrogate	Rec. (%)	Control Limits
Dibromofluoromethane	96	80-126
1,2-Dichloroethane-d4	100	80-134
Toluene-d8	99	80-120
Toluene-d8-TPPH	95	88-112
1,4-Bromofluorobenzene	97	80-120





Environmental Science International, Inc. 354 Uluniu Street, Suite 304

Kailua, HI 96734-2500

Date Received:

10/29/14 14-10-2234

Work Order: Preparation:

EPA 5030C

Method:

GC/MS / EPA 8260B

Units:

ug/L

Project: Red Hill LTM 112066

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Client Sample Number	Lab Sample Number	Date/Time Collected	Matrix	Instrument	Date Prepared	Date/Time Analyzed	QC Batch II
ES117	14-10-2234-2-A	10/28/14 12:00	Aqueous	GC/MS OO	10/31/14	11/01/14 02:36	141031L04
Comment(s): - Results were eval	luated to the MDL (DL), con	centrations >=	to the MDL (DL	but < RL (LOC	Q), if found, a	re qualified with	a "J" flag.
<u>Parameter</u>	<u>Result</u>	<u>DL</u>	<u>LOD</u>	<u>LO0</u>	<u>3</u>	<u>DF</u>	Qualifiers
Acetone	<10	6.0	10	20		1.00	U,ICH
Benzene	<0.50	0.14	0.50	1.0		1.00	U
Bromodichloromethane	<0.50	0.21	0.50	5.0		1.00	U
Bromoform	<1.0	0.50	1.0	10		1.00	U
Bromomethane	<5.0	3.9	5.0	20		1.00	U,IJ
2-Butanone	<5.0	2.2	5.0	10		1.00	U
Carbon Tetrachloride	<0.50	0.23	0.50	1.0		1.00	U
Chlorobenzene	<0.50	0.17	0.50	5.0		1.00	U
Chloroethane	<5.0	2.3	5.0	10		1.00	U
Chloroform	<0.50	0.46	0.50	5.0		1.00	U
Chloromethane	<2.0	1.8	2.0	10		1.00	U
Dibromochloromethane	<0.50	0.25	0.50	1.0		1.00	U
1,2-Dibromo-3-Chloropropane	<2.0	1.2	2.0	10		1.00	U
1,2-Dibromoethane	<0.50	0.36	0.50	1.0		1.00	U
1,2-Dichlorobenzene	<0.50	0.46	0.50	1.0		1.00	U
1,3-Dichlorobenzene	<0.50	0.40	0.50	1.0		1.00	U
1,4-Dichlorobenzene	<0.50	0.43	0.50	1.0		1.00	U
1,1-Dichloroethane	<0.50	0.28	0.50	5.0		1.00	U
1,2-Dichloroethane	<0.50	0.24	0.50	1.0		1.00	U
1,1-Dichloroethene	<0.50	0.43	0.50	1.0		1.00	U
c-1,2-Dichloroethene	<0.50	0.48	0.50	1.0		1.00	U
t-1,2-Dichloroethene	<0.50	0.37	0.50	1.0		1.00	U
1,2-Dichloropropane	<0.50	0.42	0.50	5.0		1.00	U
c-1,3-Dichloropropene	<0.50	0.25	0.50	1.0		1.00	U
t-1,3-Dichloropropene	<0.50	0.25	0.50	1.0		1.00	U
Ethylbenzene	<0.50	0.14	0.50	1.0		1.00	U
Methylene Chloride	<1.0	0.64	1.0	5.0		1.00	U
4-Methyl-2-Pentanone	<5.0	4.4	5.0	10		1.00	U
Styrene	<0.50	0.17	0.50	1.0		1.00	U
1,1,1,2-Tetrachloroethane	<0.50	0.40	0.50	1.0		1.00	U
1,1,2,2-Tetrachloroethane	<0.50	0.41	0.50	1.0		1.00	U
Tetrachloroethene	<0.50	0.39	0.50	5.0		1.00	U
Toluene	<0.50	0.24	0.50	1.0		1.00	U
1,2,4-Trichlorobenzene	<1.0	0.50	1.0	5.0		1.00	U
1,1,1-Trichloroethane	<0.50	0.30	0.50	5.0		1.00	U
Hexachloro-1,3-Butadiene	<0.50	0.32	0.50	1.0		1.00	U



Environmental Science International, Inc. 354 Uluniu Street, Suite 304 Kailua, HI 96734-2500

Work Order: Preparation: Method:

Date Received:

10/29/14 14-10-2234 **EPA 5030C**

ug/L

Units:

GC/MS / EPA 8260B

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Project: Red Hill LTM 112066

Toluene-d8

Toluene-d8-TPPH

1,4-Bromofluorobenzene

<u>Parameter</u>	Result	<u>DL</u>	<u>LOD</u>	<u>LOQ</u>	<u>DF</u>	<u>Qualifiers</u>
1,1,2-Trichloroethane	<0.50	0.38	0.50	1.0	1.00	U
Trichloroethene	<0.50	0.37	0.50	1.0	1.00	U
1,2,3-Trichloropropane	<1.0	0.64	1.0	5.0	1.00	U
Vinyl Chloride	<0.50	0.30	0.50	1.0	1.00	U
p/m-Xylene	<1.0	0.30	1.0	10	1.00	U
o-Xylene	<0.50	0.23	0.50	1.0	1.00	U
Methyl-t-Butyl Ether (MTBE)	<0.50	0.31	0.50	1.0	1.00	U
Gasoline Range Organics	<30	26	30	50	1.00	U
Surrogate	Rec. (%)	Control Limits	<u>Qualifiers</u>			
Dibromofluoromethane	96	80-126				
1,2-Dichloroethane-d4	99	80-134				

80-120

88-112

80-120

99

95

95





Environmental Science International, Inc. 354 Uluniu Street, Suite 304

Kailua, HI 96734-2500

Date Received:

10/29/14 14-10-2234

Work Order: Preparation:

EPA 5030C

Method:

GC/MS / EPA 8260B

Units:

ug/L

Project: Red Hill LTM 112066

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Client Sample Number	Lab Sample Number	Date/Time Collected	Matrix	Instrument	Date Prepared	Date/Time Analyzed	QC Batch ID
ES118	14-10-2234-3-A	10/28/14 14:30	Aqueous	GC/MS OO	10/31/14	11/01/14 03:04	141031L047
Comment(s): - Results were evaluated to	the MDL (DL), con	centrations >= t	o the MDL (DL) but < RL (LOQ), if found, are o	qualified with	a "J" flag.
<u>Parameter</u>	Result	<u>DL</u>	<u>LOD</u>	LOQ	<u>D</u> F	=	<u>Qualifiers</u>
Acetone	<10	6.0	10	20	1.0	00	U,ICH
Benzene	< 0.50	0.14	0.50	1.0	1.0	00	U
Bromodichloromethane	<0.50	0.21	0.50	5.0	1.0	00	U
Bromoform	<1.0	0.50	1.0	10	1.0	00	U
Bromomethane	<5.0	3.9	5.0	20	1.0	00	U,IJ
2-Butanone	<5.0	2.2	5.0	10	1.0	00	U
Carbon Tetrachloride	<0.50	0.23	0.50	1.0	1.0	00	U
Chlorobenzene	<0.50	0.17	0.50	5.0	1.0	00	U
Chloroethane	<5.0	2.3	5.0	10	1.0	00	U
Chloroform	<0.50	0.46	0.50	5.0	1.0	00	U
Chloromethane	<2.0	1.8	2.0	10	1.0	00	U
Dibromochloromethane	<0.50	0.25	0.50	1.0	1.0	00	U
1,2-Dibromo-3-Chloropropane	<2.0	1.2	2.0	10	1.0	00	U
1,2-Dibromoethane	<0.50	0.36	0.50	1.0	1.0	00	U
1,2-Dichlorobenzene	<0.50	0.46	0.50	1.0	1.0	00	U
1,3-Dichlorobenzene	<0.50	0.40	0.50	1.0	1.0	00	U
1,4-Dichlorobenzene	<0.50	0.43	0.50	1.0	1.0	00	U
1,1-Dichloroethane	<0.50	0.28	0.50	5.0	1.0	00	U
1,2-Dichloroethane	<0.50	0.24	0.50	1.0	1.0	00	U
1,1-Dichloroethene	<0.50	0.43	0.50	1.0	1.0	00	U
c-1,2-Dichloroethene	<0.50	0.48	0.50	1.0	1.0	00	U
t-1,2-Dichloroethene	<0.50	0.37	0.50	1.0	1.0	00	U
1,2-Dichloropropane	<0.50	0.42	0.50	5.0	1.0	00	U
c-1,3-Dichloropropene	<0.50	0.25	0.50	1.0	1.0	00	U
t-1,3-Dichloropropene	<0.50	0.25	0.50	1.0	1.0	00	U
Ethylbenzene	<0.50	0.14	0.50	1.0	1.0	00	U
Methylene Chloride	<1.0	0.64	1.0	5.0	1.0	00	U
4-Methyl-2-Pentanone	<5.0	4.4	5.0	10	1.0	00	U
Styrene	<0.50	0.17	0.50	1.0	1.0	00	U
1,1,1,2-Tetrachloroethane	<0.50	0.40	0.50	1.0	1.0		U
1,1,2,2-Tetrachloroethane	<0.50	0.41	0.50	1.0	1.0		U
Tetrachloroethene	<0.50	0.39	0.50	5.0	1.0		U
Toluene	<0.50	0.24	0.50	1.0	1.0		U
1,2,4-Trichlorobenzene	<1.0	0.50	1.0	5.0	1.0		U
1,1,1-Trichloroethane	<0.50	0.30	0.50	5.0	1.0		U
Hexachloro-1,3-Butadiene	<0.50	0.32	0.50	1.0	1.0		U



Environmental Science International, Inc. 354 Uluniu Street, Suite 304 Kailua, HI 96734-2500

Date Received: Work Order: Preparation:

<u>LOD</u>

<u>LOQ</u>

<u>DF</u>

10/29/14 14-10-2234 EPA 5030C

Method:

GC/MS / EPA 8260B

ug/L

Units:

<u>DL</u>

80-120

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Qualifiers U U U U U U U U

Project: Red H	lill LTM 112066
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<u>Parameter</u>

1,4-Bromofluorobenzene

				_
<0.50	0.38	0.50	1.0	1.00
<0.50	0.37	0.50	1.0	1.00
<1.0	0.64	1.0	5.0	1.00
<0.50	0.30	0.50	1.0	1.00
<1.0	0.30	1.0	10	1.00
<0.50	0.23	0.50	1.0	1.00
<0.50	0.31	0.50	1.0	1.00
<30	26	30	50	1.00
Rec. (%)	Control Limits	Qualifiers		
96	80-126			
100	80-134			
99	80-120			
96	88-112			
	<0.50 <1.0 <0.50 <1.0 <0.50 <0.50 <30 Rec. (%) 96 100 99	<0.50	<0.50	<0.50

Result

97





Environmental Science International, Inc.

Kailua, HI 96734-2500

Date Received: Work Order:

10/29/14 14-10-2234

354 Uluniu Street, Suite 304

Preparation:

EPA 5030C

Method:

GC/MS / EPA 8260B

Units:

ug/L

Project: Red Hill LTM 112066

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Client Sample Number	Lab Sample Number	Date/Time Collected	Matrix	Instrument	Date Prepared	Date/Time Analyzed	QC Batch ID
Method Blank	099-13-057-72	N/A	Aqueous	GC/MS OO	10/31/14	10/31/14 22:31	141031L047
Comment(s): - Results were evaluated	d to the MDL (DL), cor	ncentrations >=	to the MDL (DL) but < RL (LOC)), if found, are	qualified with	a "J" flag.
<u>Parameter</u>	Result	<u>DL</u>	LOD	LOC	<u> </u>	<u>)F</u>	Qualifiers
Acetone	<10	6.0	10	20	1	.00	U
Benzene	<0.50	0.14	0.50	1.0	1	.00	U
Bromodichloromethane	< 0.50	0.21	0.50	5.0	1	.00	U
Bromoform	<1.0	0.50	1.0	10	1	.00	U
Bromomethane	<5.0	3.9	5.0	20	1	.00	U
2-Butanone	<5.0	2.2	5.0	10	1	.00	U
Carbon Tetrachloride	<0.50	0.23	0.50	1.0	1	.00	U
Chlorobenzene	<0.50	0.17	0.50	5.0	1	.00	U
Chloroethane	<5.0	2.3	5.0	10	1	.00	U
Chloroform	<0.50	0.46	0.50	5.0	1	.00	U
Chloromethane	<2.0	1.8	2.0	10	1	.00	U
Dibromochloromethane	< 0.50	0.25	0.50	1.0	1	.00	U
1,2-Dibromo-3-Chloropropane	<2.0	1.2	2.0	10	1	.00	U
1,2-Dibromoethane	<0.50	0.36	0.50	1.0	1	.00	U
1,2-Dichlorobenzene	<0.50	0.46	0.50	1.0		.00	U
1,3-Dichlorobenzene	<0.50	0.40	0.50	1.0	1	.00	U
1,4-Dichlorobenzene	<0.50	0.43	0.50	1.0		.00	U
1,1-Dichloroethane	<0.50	0.28	0.50	5.0	1	.00	U
1,2-Dichloroethane	<0.50	0.24	0.50	1.0		.00	U
1,1-Dichloroethene	<0.50	0.43	0.50	1.0		.00	U
c-1,2-Dichloroethene	<0.50	0.48	0.50	1.0		.00	U
t-1,2-Dichloroethene	<0.50	0.37	0.50	1.0		.00	U
1,2-Dichloropropane	<0.50	0.42	0.50	5.0	1	.00	U
c-1,3-Dichloropropene	<0.50	0.25	0.50	1.0	1	.00	U
t-1,3-Dichloropropene	<0.50	0.25	0.50	1.0		.00	U
Ethylbenzene	<0.50	0.14	0.50	1.0		.00	U
Methylene Chloride	<1.0	0.64	1.0	5.0		.00	U
4-Methyl-2-Pentanone	<5.0	4.4	5.0	10		.00	U
Styrene	<0.50	0.17	0.50	1.0		.00	U
1,1,1,2-Tetrachloroethane	<0.50	0.40	0.50	1.0		.00	U
1,1,2,2-Tetrachloroethane	<0.50	0.41	0.50	1.0		.00	U
Tetrachloroethene	<0.50	0.39	0.50	5.0		.00	U
Toluene	<0.50	0.24	0.50	1.0		.00	U
1,2,4-Trichlorobenzene	<1.0	0.50	1.0	5.0		.00	U
1,1,1-Trichloroethane	<0.50	0.30	0.50	5.0		.00	U
Hexachloro-1,3-Butadiene	<0.50	0.32	0.50	1.0		.00	U



Environmental Science International, Inc. 354 Uluniu Street, Suite 304

Kailua, HI 96734-2500

1,4-Bromofluorobenzene

Date Received: Work Order:

Preparation: Method:

Units:

14-10-2234 EPA 5030C

GC/MS / EPA 8260B

ug/L

10/29/14

Project: Red Hill LTM 112066

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Doromotor	Dooult	DI	LOD	1.00	DE	Ouglifiere
<u>Parameter</u>	<u>Result</u>	<u>DL</u>	<u>LOD</u>	<u>LOQ</u>	<u>DF</u>	<u>Qualifiers</u>
1,1,2-Trichloroethane	< 0.50	0.38	0.50	1.0	1.00	U
Trichloroethene	<0.50	0.37	0.50	1.0	1.00	U
1,2,3-Trichloropropane	<1.0	0.64	1.0	5.0	1.00	U
Vinyl Chloride	<0.50	0.30	0.50	1.0	1.00	U
p/m-Xylene	<1.0	0.30	1.0	10	1.00	U
o-Xylene	<0.50	0.23	0.50	1.0	1.00	U
Methyl-t-Butyl Ether (MTBE)	<0.50	0.31	0.50	1.0	1.00	U
Gasoline Range Organics	<30	26	30	50	1.00	U
<u>Surrogate</u>	Rec. (%)	Control Limits	Qualifiers			
Dibromofluoromethane	95	80-126				
1,2-Dichloroethane-d4	99	80-134				
Toluene-d8	99	80-120				
Toluene-d8-TPPH	95	88-112				

80-120

96



Quality Control - Spike/Spike Duplicate

Environmental Science International, Inc.

Date Received:

Work Order:

10/29/14

Work Order:

14-10-2234

Kailua, HI 96734-2500

Preparation:

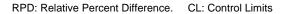
Method:

EPA 3005A Filt.

Method:

Project: Red Hill LTM 112066 Page 1 of 4

Quality Control Sample ID	Туре		Matrix	Matrix Instrument		Date Prepared Date A		lyzed	zed MS/MSD Batch Number	
ES118	Sample		Aqueous		P/MS 04	10/31/14	11/03/14	15:10	141031S04	
ES118	Matrix Spike		Aqueous ICP/		P/MS 04	10/31/14 11/03/		14:59	141031S04	
ES118	Matrix Spike Duplicate		Aqueou	Aqueous ICP/MS 04		10/31/14	11/03/14	15:01	141031S04	
Parameter	Sample Conc.	<u>Spike</u> Added	MS Conc.	MS %Rec.	MSD Conc.	MSD %Rec.	%Rec. CL	RPD	RPD CL	Qualifiers
Lead	ND	100.0	109.0	109	109.2	109	80-120	0	0-20	





Quality Control - Spike/Spike Duplicate

Environmental Science International, Inc. 354 Uluniu Street, Suite 304

Kailua, HI 96734-2500

Project: Red Hill LTM 112066

Date Received:

Work Order:

Method:

Preparation:

10/29/14 14-10-2234

EPA 3510C

EPA 8270C SIM PAHs

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Quality Control Sample ID	Туре		Matrix Instrument		Date Prepared Date Analyzed		llyzed	MS/MSD Batch Number			
14-10-2131-3	Sample	Sample		ıs (GC/MS AAA	10/30/14	10/30/14 11/03/14 18:28		141030S08		
14-10-2131-3	Matrix Spike	Matrix Spike		Aqueous GC/MS AAA		10/30/14	10/30/14 10/31/14 18:04		141030S08		
14-10-2131-3	Matrix Spike	Duplicate	Aqueous		GC/MS AAA	10/30/14	10/31/14	10/31/14 18:27		141030S08	
<u>Parameter</u>	<u>Sample</u> <u>Conc.</u>	<u>Spike</u> Added	MS Conc.	MS %Rec	MSD Conc.	MSD %Rec.	%Rec. CL	RPD	RPD CL	Qualifiers	
Naphthalene	143.3	2.000	83.89	0	78.94	0	21-133	6	0-25	3	
2-Methylnaphthalene	42.74	2.000	34.71	0	32.01	0	21-140	8	0-25	3	
1-Methylnaphthalene	59.00	2.000	47.61	0	45.00	0	20-140	6	0-25	3	
Acenaphthylene	ND	2.000	1.706	85	1.549	77	33-145	10	0-25		
Acenaphthene	0.5300	2.000	2.109	79	1.970	72	49-121	7	0-25		
Fluorene	ND	2.000	1.860	93	1.701	85	59-121	9	0-25		
Phenanthrene	ND	2.000	1.624	81	1.459	73	54-120	11	0-25		
Anthracene	ND	2.000	1.548	77	1.417	71	27-133	9	0-25		
Fluoranthene	ND	2.000	1.666	83	1.524	76	26-137	9	0-25		
Pyrene	ND	2.000	1.581	79	1.473	74	18-168	7	0-25		
Benzo (a) Anthracene	ND	2.000	1.551	78	1.422	71	33-143	9	0-25		
Chrysene	ND	2.000	1.540	77	1.402	70	17-168	9	0-25		
Benzo (k) Fluoranthene	ND	2.000	1.447	72	1.319	66	24-159	9	0-25		
Benzo (b) Fluoranthene	ND	2.000	1.372	69	1.248	62	24-159	9	0-25		
Benzo (a) Pyrene	ND	2.000	1.331	67	1.216	61	17-163	9	0-25		
Indeno (1,2,3-c,d) Pyrene	ND	2.000	1.320	66	1.166	58	10-171	12	0-25		
Dibenz (a,h) Anthracene	ND	2.000	1.422	71	1.275	64	10-219	11	0-25		
Benzo (g,h,i) Perylene	ND	2.000	1.524	76	1.373	69	10-227	10	0-25		

RPD: Relative Percent Difference. CL: Control Limits



Quality Control - Spike/Spike Duplicate

Environmental Science International, Inc.

Date Received:

Work Order:

10/29/14

Work Order:

14-10-2234

Kailua, HI 96734-2500

Preparation:

EPA 5030C

Method: GC/MS / EPA 8260B

Project: Red Hill LTM 112066 Page 3 of 4

Quality Control Sample ID	Туре	Туре			Instrument	Date Prepared	Date Ana	ate Analyzed MS/MSD Batch Nur //31/14 22:58 141031S031			
14-10-2387-2	Sample	Sample		(GC/MS OO	10/31/14	10/31/14	22:58	141031S031		
14-10-2387-2	Matrix Spike		Aqueous	(GC/MS OO	10/31/14	10/31/14	23:26	141031 S 031		
14-10-2387-2	Matrix Spike	Duplicate	Aqueous		GC/MS OO	10/31/14	10/31/14	23:53	141031 S 031		
Parameter	Sample Conc.	<u>Spike</u> Added	MS Conc.	MS %Red	MSD Conc.	MSD %Rec.	%Rec. CL	RPD	RPD CL	Qualifiers	
Acetone	ND	50.00	67.72	135	76.03	152	40-140	12	0-20	3	
Benzene	ND	50.00	49.93	100	49.48	99	80-120	1	0-20		
Bromodichloromethane	ND	50.00	48.33	97	48.57	97	75-120	0	0-20		
Bromoform	ND	50.00	49.21	98	50.44	101	70-130	2	0-20		
Bromomethane	ND	50.00	49.22	98	41.35	83	30-145	17	0-20		
2-Butanone	ND	50.00	59.57	119	59.62	119	30-150	0	0-20		
Carbon Tetrachloride	ND	50.00	51.03	102	49.80	100	65-140	2	0-20		
Chlorobenzene	ND	50.00	48.15	96	48.88	98	80-120	2	0-20		
Chloroethane	ND	50.00	44.85	90	44.26	89	60-135	1	0-20		
Chloroform	ND	50.00	50.30	101	49.66	99	65-135	1	0-20		
Chloromethane	ND	50.00	44.19	88	44.72	89	40-125	1	0-20		
Dibromochloromethane	ND	50.00	49.33	99	50.66	101	60-135	3	0-20		
1,2-Dibromo-3-Chloropropane	ND	50.00	43.24	86	46.97	94	50-130	8	0-20		
1,2-Dibromoethane	ND	50.00	50.20	100	52.08	104	80-120	4	0-20		
1,2-Dichlorobenzene	ND	50.00	47.25	95	48.04	96	70-120	2	0-20		
1,3-Dichlorobenzene	ND	50.00	48.17	96	47.98	96	75-125	0	0-20		
1,4-Dichlorobenzene	ND	50.00	46.86	94	47.50	95	75-125	1	0-20		
1,1-Dichloroethane	ND	50.00	48.12	96	47.56	95	70-135	1	0-20		
1,2-Dichloroethane	ND	50.00	49.85	100	50.61	101	70-130	1	0-20		
1,1-Dichloroethene	ND	50.00	46.89	94	46.82	94	70-130	0	0-20		
c-1,2-Dichloroethene	ND	50.00	51.13	102	50.25	101	70-125	2	0-20		
t-1,2-Dichloroethene	ND	50.00	50.60	101	49.80	100	60-140	2	0-20		
1,2-Dichloropropane	ND	50.00	48.06	96	48.21	96	75-125	0	0-20		
c-1,3-Dichloropropene	ND	50.00	51.62	103	52.11	104	70-130	1	0-20		
t-1,3-Dichloropropene	ND	50.00	50.09	100	51.77	104	55-140	3	0-20		
Ethylbenzene	ND	50.00	49.56	99	49.72	99	75-125	0	0-20		
Methylene Chloride	ND	50.00	47.52	95	49.08	98	55-140	3	0-20		
4-Methyl-2-Pentanone	ND	50.00	44.42	89	50.80	102	60-135	13	0-20		
Styrene	ND	50.00	49.18	98	49.95	100	65-135	2	0-20		
1,1,1,2-Tetrachloroethane	ND	50.00	50.43	101	50.94	102	80-130	1	0-20		
1,1,2,2-Tetrachloroethane	ND	50.00	48.99	98	51.35	103	65-130	5	0-20		
Tetrachloroethene	ND	50.00	50.71	101	52.32	105	45-150	3	0-20		
Toluene	ND	50.00	48.58	97	49.23	98	75-120	1	0-20		
1,2,4-Trichlorobenzene	ND	50.00	44.33	89	46.15	92	65-135	4	0-20		
1,1,1-Trichloroethane	ND	50.00	50.20	100	48.87	98	65-130	3	0-20		



Quality Control - Spike/Spike Duplicate

Environmental Science International, Inc.

Date Received:

Work Order:

10/29/14

Work Order:

14-10-2234

Kailua, HI 96734-2500

Preparation:

Method:

GC/MS / EPA 8260B

Project: Red Hill LTM 112066 Page 4 of 4

<u>Parameter</u>	Sample Conc.	<u>Spike</u> Added	MS Conc.	MS %Rec.	MSD Conc.	MSD %Rec.	%Rec. CL	RPD	RPD CL	<u>Qualifiers</u>
Hexachloro-1,3-Butadiene	ND	50.00	45.69	91	45.79	92	50-140	0	0-20	
1,1,2-Trichloroethane	ND	50.00	49.59	99	51.28	103	75-125	3	0-20	
Trichloroethene	ND	50.00	47.79	96	47.66	95	70-125	0	0-20	
1,2,3-Trichloropropane	ND	50.00	47.54	95	51.06	102	75-125	7	0-20	
Vinyl Chloride	ND	50.00	47.42	95	46.59	93	50-145	2	0-20	
p/m-Xylene	ND	100.0	98.13	98	98.70	99	75-130	1	0-20	
o-Xylene	ND	50.00	49.82	100	49.94	100	80-120	0	0-20	
Methyl-t-Butyl Ether (MTBE)	ND	50.00	47.63	95	49.54	99	65-125	4	0-20	



Quality Control - PDS

Environmental Science International, Inc.

Date Received:

Work Order:

10/29/14

Work Order:

14-10-2234

Kailua, HI 96734-2500

Preparation:

Method:

EPA 3005A Filt.

Method:

Quality Control Sample ID	Туре	N	Matrix	Instrument	Date Prepared	Date Analyzed	PDS/PDSD Batch Number
ES118	Sample	A	Aqueous	ICP/MS 04	10/31/14 00:00	11/03/14 15:10	141031S04
ES118	PDS	A	Aqueous	ICP/MS 04	10/31/14 00:00	11/03/14 15:03	141031S04
<u>Parameter</u>		Sample Conc.	Spike Added	PDS Conc.	PDS %Re	ec. %Rec. C	<u>Qualifiers</u>
Lead		ND	100.0	102.6	103	75-125	





Quality Control - LCS/LCSD

Environmental Science International, Inc.

Date Received:

Work Order:

10/29/14

Work Order:

14-10-2234

Kailua, HI 96734-2500

Preparation:

EPA 8015B (M)

Quality Control Sample ID	Туре	Matrix	Instrument	Date Prepa	ared Date	Analyzed	LCS/LCSD Ba	atch Number
099-15-516-206	LCS	Aqueou	s GC 45	10/30/14	10/31	1/14 07:20	141030B06	
099-15-516-206	LCSD	Aqueou	s GC 45	10/30/14	10/31	1/14 07:37	141030B06	
Parameter	Spike Added LCS	Conc. LC %F		LCSD 9	%Rec. CL	RPD	RPD CL	Qualifiers
TPH as Diesel	2000 229	6 11	2219	111	60-132	3	0-11	



Quality Control - LCS

Environmental Science International, Inc.

Date Received:

Work Order:

10/29/14

Work Order:

14-10-2234

Kailua, HI 96734-2500

Preparation:

Method:

EPA 3005A Filt.

Method:

Project: Red Hill LTM 112066 Page 2 of 5

Quality Control Sample ID	Туре	Matrix Instrument Date Prepare		Date Prepared	Date Analyzed	LCS Batch Number
099-14-497-104	LCS	Aqueous	ICP/MS 04	10/31/14	11/03/14 14:50	141031L04F
<u>Parameter</u>		Spike Added	Conc. Recovere	ed LCS %Re	ec. %Rec	. CL Qualifiers
Lead		100.0	102.6	103	80-12	0



Quality Control - LCS

Environmental Science International, Inc. 354 Uluniu Street, Suite 304

Kailua, HI 96734-2500

Project: Red Hill LTM 112066

Date Received:

14-10-2234

10/29/14

Work Order: Preparation:

EPA 3510C

Method:

EPA 8270C SIM PAHs

Page 3 of 5

Quality Control Sample ID	Туре	Matrix	Instrument	Date Prepared D	Date Analyzed LCS Ba	atch Number
099-15-148-67	LCS	Aqueous	GC/MS AAA	10/30/14 1	0/31/14 17:40 141030	DL08
Parameter		Spike Added	Conc. Recove	ered LCS %Rec.	%Rec. CL	Qualifiers
Naphthalene		2.000	1.394	70	21-133	
2-Methylnaphthalene		2.000	1.382	69	21-140	
1-Methylnaphthalene		2.000	1.409	70	20-140	
Acenaphthylene		2.000	1.394	70	33-145	
Acenaphthene		2.000	1.451	73	55-121	
Fluorene		2.000	1.539	77	59-121	
Phenanthrene		2.000	1.447	72	54-120	
Anthracene		2.000	1.308	65	27-133	
Fluoranthene		2.000	1.538	77	26-137	
Pyrene		2.000	1.480	74	45-129	
Benzo (a) Anthracene		2.000	1.398	70	33-143	
Chrysene		2.000	1.425	71	17-168	
Benzo (k) Fluoranthene		2.000	1.412	71	24-159	
Benzo (b) Fluoranthene		2.000	1.322	66	24-159	
Benzo (a) Pyrene		2.000	1.150	58	17-163	
Indeno (1,2,3-c,d) Pyrene		2.000	1.251	63	25-175	
Dibenz (a,h) Anthracene		2.000	1.391	70	25-175	
Benzo (g,h,i) Perylene		2.000	1.514	76	25-157	



Quality Control - LCS/LCSD

Environmental Science International, Inc. 354 Uluniu Street, Suite 304

Kailua, HI 96734-2500

Date Received:

Method:

Work Order:
Preparation:

GC/MS / EPA 8260B

10/29/14

14-10-2234

EPA 5030C

Project: Red Hill LTM 112066 Page 4 of 5

Quality Control Sample ID	Туре	Mat	trix	Instrument	Date Pre	pared	Date	Analyzed	LCS/LCSD Ba	atch Number
099-13-057-72	LCS	Aqı	ueous	GC/MS OO	10/31/14		10/31	1/14 21:09	141031L047	
099-13-057-72	LCSD	Aqı	ueous	GC/MS OO	10/31/14		10/31	/14 21:37	141031L047	
Parameter	Spike Added	LCS Conc.	<u>LCS</u> %Rec.	LCSD Conc.	LCSD %Rec.	%Red	:. CL	RPD	RPD CL	Qualifiers
Acetone	50.00	66.09	132	N/A	N/A	40-14	0	N/A	0-20	
Benzene	50.00	50.90	102	N/A	N/A	80-12	0	N/A	0-20	
Bromodichloromethane	50.00	49.29	99	N/A	N/A	75-12	0	N/A	0-20	
Bromoform	50.00	50.67	101	N/A	N/A	70-13	0	N/A	0-20	
Bromomethane	50.00	43.60	87	N/A	N/A	30-14	5	N/A	0-20	
2-Butanone	50.00	55.76	112	N/A	N/A	30-15	0	N/A	0-20	
Carbon Tetrachloride	50.00	52.82	106	N/A	N/A	65-14	0	N/A	0-20	
Chlorobenzene	50.00	50.31	101	N/A	N/A	80-12	0	N/A	0-20	
Chloroethane	50.00	46.45	93	N/A	N/A	60-13	5	N/A	0-20	
Chloroform	50.00	51.26	103	N/A	N/A	65-13	5	N/A	0-20	
Chloromethane	50.00	46.27	93	N/A	N/A	40-12	5	N/A	0-20	
Dibromochloromethane	50.00	50.80	102	N/A	N/A	60-13	5	N/A	0-20	
1,2-Dibromo-3-Chloropropane	50.00	47.64	95	N/A	N/A	50-13	0	N/A	0-20	
1,2-Dibromoethane	50.00	52.19	104	N/A	N/A	80-12	0	N/A	0-20	
1,2-Dichlorobenzene	50.00	49.11	98	N/A	N/A	70-12	0	N/A	0-20	
1,3-Dichlorobenzene	50.00	49.80	100	N/A	N/A 75-125		25 N/A		0-20	
1,4-Dichlorobenzene	50.00	49.36	99	N/A	N/A	75-12	5	N/A	0-20	
1,1-Dichloroethane	50.00	50.29	101	N/A	N/A	70-13	5	N/A	0-20	
1,2-Dichloroethane	50.00	51.26	103	N/A	N/A	70-13	0	N/A	0-20	
1,1-Dichloroethene	50.00	50.07	100	N/A	N/A	70-130 N/		N/A	0-20	
c-1,2-Dichloroethene	50.00	52.18	104	N/A	N/A	70-12	5	N/A	0-20	
t-1,2-Dichloroethene	50.00	53.26	107	N/A	N/A	60-14	0	N/A	0-20	
1,2-Dichloropropane	50.00	49.39	99	N/A	N/A	75-12	5	N/A	0-20	
c-1,3-Dichloropropene	50.00	53.06	106	N/A	N/A	70-13	0	N/A	0-20	
t-1,3-Dichloropropene	50.00	53.91	108	N/A	N/A	55-14	0	N/A	0-20	
Ethylbenzene	50.00	52.31	105	N/A	N/A	75-125 N/A 0-20				
Methylene Chloride	50.00	50.23	100	N/A	N/A	55-14	0	N/A	0-20	
4-Methyl-2-Pentanone	50.00	49.85	100	N/A	N/A	60-13	5	N/A	0-20	
Styrene	50.00	51.56	103	N/A	N/A	65-13	5	N/A	0-20	
1,1,1,2-Tetrachloroethane	50.00	51.98	104	N/A	N/A	80-13	0	N/A	0-20	
1,1,2,2-Tetrachloroethane	50.00	50.95	102	N/A	N/A	65-13	0	N/A	0-20	
Tetrachloroethene	50.00	51.34	103	N/A	N/A	45-15	0	N/A	0-20	
Toluene	50.00	50.99	102	N/A	N/A	75-12	0	N/A	0-20	
1,2,4-Trichlorobenzene	50.00	47.93	96	N/A	N/A	65-13	5	N/A	0-20	
1,1,1-Trichloroethane	50.00	52.27	105	N/A	N/A	65-13	0	N/A	0-20	
Hexachloro-1,3-Butadiene	50.00	49.37	99	N/A	N/A	50-14	0	N/A	0-20	





Quality Control - LCS/LCSD

Environmental Science International, Inc.

Date Received: 10/29/14

Work Order: 14-10-2234

Kailua, HI 96734-2500 Preparation: EPA 5030C

Method: GC/MS / EPA 8260B

Project: Red Hill LTM 112066 Page 5 of 5

<u>Parameter</u>	Spike Added	LCS Conc.	LCS %Rec.	LCSD Conc.	LCSD %Rec.	%Rec. CL	<u>RPD</u>	RPD CL	<u>Qualifiers</u>
1,1,2-Trichloroethane	50.00	50.61	101	N/A	N/A	75-125	N/A	0-20	
Trichloroethene	50.00	50.27	101	N/A	N/A	70-125	N/A	0-20	
1,2,3-Trichloropropane	50.00	50.55	101	N/A	N/A	75-125	N/A	0-20	
Vinyl Chloride	50.00	49.26	99	N/A	N/A	50-145	N/A	0-20	
p/m-Xylene	100.0	103.2	103	N/A	N/A	75-130	N/A	0-20	
o-Xylene	50.00	51.96	104	N/A	N/A	80-120	N/A	0-20	
Methyl-t-Butyl Ether (MTBE)	50.00	50.09	100	N/A	N/A	65-125	N/A	0-20	
Gasoline Range Organics	1000	1009	101	1025	103	80-120	2	0-20	



Sample Analysis Summary Report

Work Order: 14-10-2234				Page 1 of 1
Method	Extraction	Chemist ID	Instrument	Analytical Location
EPA 6020	EPA 3005A Filt.	598	ICP/MS 04	1
EPA 8015B (M)	EPA 3510C	682	GC 45	1
EPA 8270C SIM PAHs	EPA 3510C	923	GC/MS AAA	1
GC/MS / EPA 8260B	EPA 5030C	849	GC/MS OO	2

Location 1: 7440 Lincoln Way, Garden Grove, CA 92841 Location 2: 7445 Lampson Avenue, Garden Grove, CA 92841



Glossary of Terms and Qualifiers

Work Order: 14-10-2234 Page 1 of 1

Qualifiers	<u>Definition</u>
*	See applicable analysis comment.
<	Less than the indicated value.
>	Greater than the indicated value.
1	Surrogate compound recovery was out of control due to a required sample dilution. Therefore, the sample data was reported without further clarification.
2	Surrogate compound recovery was out of control due to matrix interference. The associated method blank surrogate spike compound was in control and, therefore, the sample data was reported without further clarification.
3	Recovery of the Matrix Spike (MS) or Matrix Spike Duplicate (MSD) compound was out of control due to suspected matrix interference. The associated LCS recovery was in control.
4	The MS/MSD RPD was out of control due to suspected matrix interference.
5	The PDS/PDSD or PES/PESD associated with this batch of samples was out of control due to suspected matrix interference.
6	Surrogate recovery below the acceptance limit.
7	Surrogate recovery above the acceptance limit.
В	Analyte was present in the associated method blank.
BU	Sample analyzed after holding time expired.
BV	Sample received after holding time expired.
DL	The Detection Limit (DL) is the smallest analyte concentration that can be demonstrated to be different from zero or a blank concentration at the 99% level of confidence.
E	Concentration exceeds the calibration range.
ET	Sample was extracted past end of recommended max. holding time.
HD	The chromatographic pattern was inconsistent with the profile of the reference fuel standard.
ICH	Initial calibration verification recovery is above the control limit for this analyte.
ICJ	Initial calibration verification recovery is below the control limit for this analyte.
IH	Calibration verification recovery is above the control limit for this analyte.
IJ	Calibration verification recovery is below the control limit for this analyte.
J	Analyte was detected at a concentration below the LOQ and above the DL. Reported value is estimated.
JA	Analyte positively identified but quantitation is an estimate.
LOD	The Limit of Detection (LOD) is the smallest amount or concentration of a substance that must be present in a sample in order to be detected at 99% confidence level.
LOQ	The Limit of Quantitation (LOQ) is the lowest concentration of a substance that produces a quantitative result within specified limits of precision and bias.
Q	Spike recovery and RPD control limits do not apply resulting from the parameter concentration in the sample exceeding the spike concentration by a factor of four or greater.

- SG The sample extract was subjected to Silica Gel treatment prior to analysis.
- U Undetected at Detection Limit (DL) and is reported as less than the Limit of Detection (LOD).
- X % Recovery and/or RPD out-of-range.
- Z Analyte presence was not confirmed by second column or GC/MS analysis.

Solid - Unless otherwise indicated, solid sample data is reported on a wet weight basis, not corrected for % moisture. All QC results are reported on a wet weight basis.

Any parameter identified in 40CFR Part 136.3 Table II that is designated as "analyze immediately" with a holding time of <= 15 minutes (40CFR-136.3 Table II, footnote 4), is considered a "field" test and the reported results will be qualified as being received outside of the stated holding time unless received at the laboratory within 15 minutes of the collection time.

A calculated total result (Example: Total Pesticides) is the summation of each component concentration and/or, if "J" flags are reported, estimated concentration. Component concentrations showing not detected (ND) are summed into the calculated total result as zero concentrations.

Calscience Environmental Laboratories, Inc.

CHAIN OF CUSTODY RECORD

Date 10/28/14

WO #/LAB USE ONLY

7440 Lincoln Way, Garden Grove, CA 92841-1427 • (714) 895-5494 Other locations: Concord, San Luis Obispo, Houston, and Corpus Christi For courier service / sample drop off information,

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DISTRIBUTION: White with final report, Green and Yellow to Client.

Please note that pages 1 and 2 of 2 of our T/Cs are printed on the reverse side of the Green and Yellow copies respectively.

ORIGIN ID:HNLA

SHIP DATE: 280CT14 ACTWGT: 52.5 LB CAD: /POS1525 DIMS: 26x15x14 IN

BILL RECIPIENT

ÚNITED STATES US

-T0

CALSCIENCE 7440 LINCOLN WAY

GARDEŃ GROVE CA 92841

(714) 895 - 5494

REF:

DEDI:



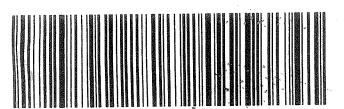
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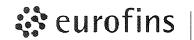
WZ APVA

92841 ca-us SNA









Calscience

SAMPLE RECEIPT FORM Cooler / of /

CLIENT: ESI	DATE:_	10/29	/ 14
TEMPERATURE: Thermometer ID: SC2 (Criteria: 0.0 °C – 6.0 °C, not froze	en except se	diment/tissu	e)
Temperature 3.5°C - 0.2°C (CF) = 3.3°C	☑Blank	☐ Sample	Э
☐ Sample(s) outside temperature criteria (PM/APM contacted by:)			
☐ Sample(s) outside temperature criteria but received on ice/chilled on same	day of sampl	ing.	
\square Received at ambient temperature, placed on ice for transport by C	ourier.		
Ambient Temperature: Ambient Temperature: Air Filter		Checked b	y: <i>[5</i>
CUSTODY SEALS INTACT:			15
☐ Cooler ☐ ☐ No (Not Intact) ☐ Not Present	□ N/A		and the
✓ Sample □ □ □ No (Not Intact) □ Not Present		Checked by	1: <u>836</u>
SAMPLE CONDITION:	Yes	No	N/A
Chain-Of-Custody (COC) document(s) received with samples			
COC document(s) received complete			
☐ Collection date/time, matrix, and/or # of containers logged in based on sample labels			
☐ No analysis requested. ☐ Not relinquished. ☐ No date/time relinquished.			
Sampler's name indicated on COC	. 🖊		
Sample container label(s) consistent with COC			
Sample container(s) intact and good condition	. 🗸		
Proper containers and sufficient volume for analyses requested	. 🗹		
Analyses received within holding time	. 🎾		
Aqueous samples received within 15-minute holding time			
☐ pH ☐ Residual Chlorine ☐ Dissolved Sulfides ☐ Dissolved Oxygen	🗆		2
Proper preservation noted on COC or sample container	,		
Unpreserved vials received for Volatiles analysis			
Volatile analysis container(s) free of headspace			
Tedlar bag(s) free of condensation CONTAINER TYPE:	🗆		4
Solid: □4ozCGJ □8ozCGJ □16ozCGJ □Sleeve () □EnCore	es [®] □Terra	Cores [®] □_	
Aqueous: ØVOA ØVOAh □VOAna₂ □125AGB □125AGBh □125AGBp	a Z1AGB [⊒1AGB na ₂ □]1AGB s
□500AGB Ø500AGJ □500AGJs □250AGB □250CGB □250CGB	s & 1PBn	□1PB na □	500PB
□250PB □250PB n □125PB □125PB znna □100PJ □100PJ na ₂ □_			
Air: □Tedlar [®] □Canister Other: □ Trip Blank Lot# : 14⊃9○8			
Container: C: Clear A: Amber P: Plastic G: Glass J: Jar B: Bottle Z: Ziploc/Resealable Bag E: E Preservative: h: HCL n: HNO ₃ na ₂ :Na ₂ S ₂ O ₃ na: NaOH p: H ₃ PO ₄ s: H ₂ SO ₄ u: Ultra-pure znna: ZnAc ₂ +N	•	Reviewed by: Scanned by:	
Fiebervative. II. FIOL II. HIVO3 Haz., Va20203 Ha. 1VaOH p. H3FO4 b. H2OO4 u. Olla-pule ziilla. ZhA02+W	Some of the second		





Supplemental Report 1

Calscience



WORK ORDER NUMBER: 14-10-2234

The difference is service



AIR | SOIL | WATER | MARINE CHEMISTRY

Analytical Report For

Client: Environmental Science International, Inc.

Client Project Name: Red Hill LTM 112066

Attention: Jeff Hattemer

354 Uluniu Street, Suite 304 Kailua, HI 96734-2500

Richard Vellas

Approved for release on 11/05/2014 by:

Richard Villafania Project Manager



ResultLink > Email your PM >

Eurofins Calscience, Inc. (Calscience) certifies that the test results provided in this report meet all NELAC requirements for parameters for which accreditation is required or available. Any exceptions to NELAC requirements are noted in the case narrative. The original report of subcontracted analyses, if any, is attached to this report. The results in this report are limited to the sample(s) tested and any reproduction thereof must be made in its entirety. The client or recipient of this report is specifically prohibited from making material changes to said report and, to the extent that such changes are made, Calscience is not responsible, legally or otherwise. The client or recipient agrees to indemnify Calscience for any defense to any litigation which may arise.



Contents

Client Project Name:	Red Hill LTM 112066
Work Order Number:	14-10-2234

1	Work Order Narrative	3
2	Client Sample Data	4
3	Quality Control Sample Data. 3.1 MS/MSD. 3.2 PDS/PDSD. 3.3 LCS/LCSD.	5
4	Sample Analysis Summary	8
5	Glossary of Terms and Qualifiers	9
6	Chain-of-Custody/Sample Receipt Form	10



Work Order Narrative

Work Order: 14-10-2234 Page 1 of 1

Condition Upon Receipt:

Samples were received under Chain-of-Custody (COC) on 10/29/14. They were assigned to Work Order 14-10-2234.

Unless otherwise noted on the Sample Receiving forms all samples were received in good condition and within the recommended EPA temperature criteria for the methods noted on the COC. The COC and Sample Receiving Documents are integral elements of the analytical report and are presented at the back of the report.

Holding Times:

All samples were analyzed within prescribed holding times (HT) and/or in accordance with the Calscience Sample Acceptance Policy unless otherwise noted in the analytical report and/or comprehensive case narrative, if required.

Any parameter identified in 40CFR Part 136.3 Table II that is designated as "analyze immediately" with a holding time of <= 15 minutes (40CFR-136.3 Table II, footnote 4), is considered a "field" test and the reported results will be qualified as being received outside of the stated holding time unless received at the laboratory within 15 minutes of the collection time.

Quality Control:

All quality control parameters (QC) were within established control limits except where noted in the QC summary forms or described further within this report.

Additional Comments:

Air - Sorbent-extracted air methods (EPA TO-4A, EPA TO-10, EPA TO-13A, EPA TO-17): Analytical results are converted from mass/sample basis to mass/volume basis using client-supplied air volumes.

New York NELAP air certification does not certify for all reported methods and analytes, reference the accredited items here: http://www.calscience.com/PDF/New_York.pdf

Solid - Unless otherwise indicated, solid sample data is reported on a wet weight basis, not corrected for % moisture. All QC results are always reported on a wet weight basis.

Subcontractor Information:

Unless otherwise noted below (or on the subcontract form), no samples were subcontracted.



Analytical Report

Environmental Science International, Inc. Date Received: 10/29/14 354 Uluniu Street, Suite 304 Work Order: 14-10-2234 Kailua, HI 96734-2500 Preparation: N/A Method: EPA 200.8 Units: ug/L

Project: Red Hill LTM 112066 Page 1 of 1

Client Sample Nu	mber	Lab Sample Number	Date/Time Collected	Matrix	Instrument	Date Prepared	Date/Time Analyzed	QC Batch ID
E3117		14-10-2234-2-J	12:00	Aqueous	ICP/IVIS U4	10/31/14	14:46	141031L01D

Comment(s): - Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are qualified with a "J" flag. <u>Parameter</u> Result <u>RL</u> <u>MDL</u> DF Qualifiers 0.211 1.00 0.0898 1.00 J Lead

Method Blank	099-16-094-581	N/A	Aqueous	ICP/MS 04	10/31/14	11/03/14 14:31	141031L01D
Comment(s):	- Results were evaluated to the MDL (DL), co	ncentrations >=	to the MDL (DL	.) but < RL (LO	Q), if found, are	qualified with a	"J" flag.
<u>Parameter</u>	Res	<u>sult</u>	<u>RL</u>	<u>MDL</u>	<u>DF</u>	<u>Q</u>	<u>ualifiers</u>
Lead	<0.	0898	1.00	0.0898	1.00	U	



RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Quality Control - Spike/Spike Duplicate

Environmental Science International, Inc.

Date Received:

Work Order:

10/29/14

Work Order:

14-10-2234

Kailua, HI 96734-2500

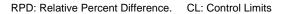
Preparation:

N/A

Method:

EPA 200.8

Quality Control Sample ID	Туре		Matrix	In	strument	Date Prepared	Date Ana	lyzed	MS/MSD Bat	ch Number
ES117	Sample		Aqueous	i IC	CP/MS 04	10/31/14	11/03/14	14:46	141031S01A	
ES117	Matrix Spike		Aqueous	s IC	CP/MS 04	10/31/14	11/03/14	14:35	141031S01A	
ES117	Matrix Spike D	uplicate	Aqueous	i IC	CP/MS 04	10/31/14	11/03/14	14:37	141031S01A	
Parameter	Sample Conc.	<u>Spike</u> <u>Added</u>	MS Conc.	MS %Rec.	MSD Conc.	MSD %Rec.	%Rec. CL	RPD	RPD CL	Qualifiers
Lead	ND	100.0	100.7	101	100.8	101	80-120	0	0-20	





Quality Control - PDS

Environmental Science International, Inc.

Date Received:

Work Order:

14-10-2234

Kailua, HI 96734-2500

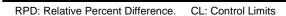
Preparation:

N/A

Method:

EPA 200.8

Quality Control Sample ID	Туре	N	Matrix	Instrument	Date Prepared	Date Analyzed	PDS/PDSD Batch Number
ES117	Sample	A	Aqueous	ICP/MS 04	10/31/14 00:00	11/03/14 14:46	141031S01A
ES117	PDS	A	Aqueous	ICP/MS 04	10/31/14 00:00	11/03/14 14:39	141031S01A
<u>Parameter</u>		Sample Conc.	Spike Added	DDS Conc	PDS %Re	ec. %Rec. C	<u>CL</u> <u>Qualifiers</u>
Lead		ND	100.0	99.82	100	75-125	





Quality Control - LCS

Environmental Science International, Inc.

Date Received:

Work Order:

14-10-2234

Kailua, HI 96734-2500

Preparation:

N/A

Method:

EPA 200.8

Quality Control Sample ID	Туре	Matrix	Instrument	Date Prepared	Date Analyzed	LCS Batch Number
099-16-094-581	LCS	Aqueous	ICP/MS 04	10/31/14	11/03/14 14:33	141031L01D
Parameter		Spike Added	Conc. Recover	red LCS %Re	ec. %Rec	. CL Qualifiers
Lead		100.0	95.21	95	80-12	0



Sample Analysis Summary Report

Work Order: 14-10-2234				Page 1 of 1
<u>Method</u>	Extraction	Chemist ID	<u>Instrument</u>	Analytical Location
EPA 200.8	N/A	598	ICP/MS 04	1



Glossary of Terms and Qualifiers

Work Order: 14-10-2234 Page 1 of 1

Qualifiers	<u>Definition</u>
*	See applicable analysis comment.
<	Less than the indicated value.
>	Greater than the indicated value.
1	Surrogate compound recovery was out of control due to a required sample dilution. Therefore, the sample data was reported without further clarification.
2	Surrogate compound recovery was out of control due to matrix interference. The associated method blank surrogate spike compound was in control and, therefore, the sample data was reported without further clarification.
3	Recovery of the Matrix Spike (MS) or Matrix Spike Duplicate (MSD) compound was out of control due to suspected matrix interference. The associated LCS recovery was in control.
4	The MS/MSD RPD was out of control due to suspected matrix interference.
5	The PDS/PDSD or PES/PESD associated with this batch of samples was out of control due to suspected matrix interference.
6	Surrogate recovery below the acceptance limit.
7	Surrogate recovery above the acceptance limit.
В	Analyte was present in the associated method blank.
BU	Sample analyzed after holding time expired.
BV	Sample received after holding time expired.
E	Concentration exceeds the calibration range.
ET	Sample was extracted past end of recommended max. holding time.
HD	The chromatographic pattern was inconsistent with the profile of the reference fuel standard.
HDH	The sample chromatographic pattern for TPH matches the chromatographic pattern of the specified standard but heavier hydrocarbons were also present (or detected).
HDL	The sample chromatographic pattern for TPH matches the chromatographic pattern of the specified standard but lighter hydrocarbons were also present (or detected).
J	Analyte was detected at a concentration below the reporting limit and above the laboratory method detection limit. Reported value is estimated.
JA	Analyte positively identified but quantitation is an estimate.
ME	LCS Recovery Percentage is within Marginal Exceedance (ME) Control Limit range (+/- 4 SD from the mean).
ND	Parameter not detected at the indicated reporting limit.
Q	Spike recovery and RPD control limits do not apply resulting from the parameter concentration in the sample exceeding the spike concentration by a factor of four or greater.
SG	The sample extract was subjected to Silica Gel treatment prior to analysis.

X % Recovery and/or RPD out-of-range.

Analyte presence was not confirmed by second column or GC/MS analysis.

Undetected at the laboratory method detection limit.

U

Ζ

Solid - Unless otherwise indicated, solid sample data is reported on a wet weight basis, not corrected for % moisture. All QC results are reported on a wet weight basis.

Any parameter identified in 40CFR Part 136.3 Table II that is designated as "analyze immediately" with a holding time of <= 15 minutes (40CFR-136.3 Table II, footnote 4), is considered a "field" test and the reported results will be qualified as being received outside of the stated holding time unless received at the laboratory within 15 minutes of the collection time.

A calculated total result (Example: Total Pesticides) is the summation of each component concentration and/or, if "J" flags are reported, estimated concentration. Component concentrations showing not detected (ND) are summed into the calculated total result as zero concentrations.

CHAIN OF CUSTODY RECORD Time: (8'002 DOD 1000 SAMPLER(S): (PRINT) oť ANALYSES X747\0208 □ X747\0108 □ slabM S27 P.O. NO. MIS 0728 X 0728 □ 2HA9 Please check box or fill in blank as needed CB2 (808S) 0 Pesticides (8081) REQUESTED Date Page SVOCs (8270) Ped HIII LTM 1120 VB Prep (5035)

Terra Core

Terra Core Oxygenates (8260) Lest lational CLIENT PROJECT NAME / NUMBER: X 14-10-2234 AOC² (85e0) □ 0928 □ 381M \ X318 WO # / LAB USE ONLY PROJECT CONTACT TPH □ C6-C36 □ C6-C44 Received by: (Signature/Affiliation) Received by: (Signature/Affiliation) (5/18) (PMS) (g)H9T**X** E-MAIL: JAMHEMERCEECHENCELORM, REPERCESHENCE OF IN Field Filtered Calscience Environmental Laboratories, Inc. LOG CODE Preserved ZIP: Other locations: Concord, San Luis Obispo, Houston, and Corpus Christi Unpreserved STATE: 96734 7440 Lincoln Way, Garden Grove, CA 92841-1427 • (714) 895-5494 STANDARD NO. OF CONT. pegular (full) Giv monitaring analytelist For courier service / sample drop off information, To S Sec. Sec. A STON MATRIX contact sales@calscience.com or call us. TURNAROUND TIME (Rush surcharges may apply to any TAT not "STANDARD") 5 DAYS I 7000 000 000 TIME SAMPLING 41/82/01 72 HR DATE ADDRESS: 354 Ulunin St 48 HR んとよう GLOBAL ID LABORATORY CLIENT: 65 SAMPLE 1D OHEO-197 (808) Relinquished by: (Signature) Relinquished by: (Signature) ☐ 24 HR SPECIAL INSTRUCTIONS: OITY: Kailya COELT EDF SAME DAY LAB USE ONLY N

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Page 10 of 12

01/01/14 Revision

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SHIP DATE: 280CT14 ACTWGT: 52.5 LB CAD: /POS1525 DIMS: 26x15x14 IN

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TO

CALSCIENCE 7440 LINCOLN WAY

GARDEN GROVE CA 92841

(714) 895-5494

REF:

NFDT:



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TRK# 8045 5791 7375

WED - 29 OCT AA STANDARD OVERNIGHT

WZ APVA

92841 ca-us SNA









Calscience

WORK ORDER #: 14-10-22

SAMPLE RECEIPT FORM Cooler / of /

CLIENT: ESI	DATE: _	10/29/	/ 14
TEMPERATURE: Thermometer ID: SC2 (Criteria: 0.0 °C − 6.0 °C, not frozen Temperature 3.5 °C − 0.2 °C (CF) = 3.3 °C Sample(s) outside temperature criteria (PM/APM contacted by:) Sample(s) outside temperature criteria but received on ice/chilled on same da Received at ambient temperature, placed on ice for transport by Cou	Blank y of sampli	☐ Sample	
Ambient Temperature: Air Filter		Checked by	y: <u> </u> 5
CUSTODY SEALS INTACT: Cooler	□ N/A	Checked by Checked by	
SAMPLE CONDITION:	'es	No	N/A
Chain-Of-Custody (COC) document(s) received with samples			
COC document(s) received complete			
☐ Collection date/time, matrix, and/or # of containers logged in based on sample labels.			
☐ No analysis requested. ☐ Not relinquished. ☐ No date/time relinquished.			
Sampler's name indicated on COC			
Sample container label(s) consistent with COC			
Sample container(s) intact and good condition			
Proper containers and sufficient volume for analyses requested	9		
Analyses received within holding time			
Aqueous samples received within 15-minute holding time			
☐ pH ☐ Residual Chlorine ☐ Dissolved Sulfides ☐ Dissolved Oxygen			
Proper preservation noted on COC or sample container			
Unpreserved vials received for Volatiles analysis			
Volatile analysis container(s) free of headspace			
Tedlar bag(s) free of condensation CONTAINER TYPE:			4
Solid: □4ozCGJ □8ozCGJ □16ozCGJ □Sleeve() □EnCores	® □Terra	Cores [®] □_	
Aqueous: ØVOA ØVOAh □VOAna₂ □125AGB □125AGBh □125AGBp [ZIAGB []1AGB na ₂ [1AGB s
□500AGB Ø500AGJ □500AGJs □250AGB □250CGB □250CGBs	Ø1PBn	□1PB na □	500PB
□250PB □250PB n □125PB □125PB znna □100PJ □100PJ na ₂ □	□		
Air: Tedlar [®] Canister Other: Trip Blank Lot#: 140908B Container: C: Clear A: Amber P: Plastic G: Glass J: Jar B: Bottle Z: Ziploc/Resealable Bag E: Enveronment	elope F	Reviewed by:	82

RAW DATA SHEET FOR METHOD: EPA 8015B (M)

WORK ORDER: **INSTRUMENT:**

14-10-2234

GC 45

D/T ANALYZED:

ANALYZED BY: 682

2014-10-31 11:16

EXTRACTION:

EPA 3510C

REVIEWED BY: D/T REVIEWED:

DATA FILE:

D/T EXTRACTED: 2014-10-30 00:00

W:\GC 45\GC 45 DATA\2014\141030\14103067.D\14103067

<u>#</u> 2

CLIENT SAMPLE NUMBER: ES117

LCS/MB BATCH:

141030B06

SAMPLE VOLUME / WEIGHT:

DEFAULT: 500.00 ml / ACTUAL: 500.00 ml

MS/MSD BATCH:

FINAL VOLUME / WEIGHT:

DEFAULT: 5.00 ml / ACTUAL: 2.50 ml

UNITS:

ug/L

ADJUSTMENT RATIO TO PF:

0.50

COMMENT:

Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, as

qualified with a "J" flag. TPH as Diesel is quantified in the carbon range C10-C28.

COMPOUND TPH as Diesel

INI. CONC DF

CONC

DL

LOD

LOQ

QUAL bJ

4440 1.00 22.2 11 12 25

Area Percent Report

Data File Name : W:\GC_45\GC 45 DATA\2014\141030\14103067.D

Page Number : 1

Operator : 682 Vial Number : Vial 67

Instrument : GC 45 Injection Number : 1 Sample Name : 14-10-2234-2 Sequence Line : 67

Instrument Method: C:\CHEM32\1\METHODS\ ->

Acquired on : 31 Oct 14 11:16 am

Report Created on: 04 Nov 14 02:51 pm Analysis Method : 8015B.MTH

Software Revision: Rev. B.03.02 [341] Copyright © Agilent Technologies

Sig. 1 in W:\GC_45\GC 45 DATA\2014\141030\ ->

Pk	Ret Time	Area	Height		Width	Response %
1						
1	2.455	0.77	0	VV	0.054	0.105
2	2.637	1.41	0	VV	0.039	0.193
3	2.694	1.80	1	VV	0.051	0.247
4	2.750	2.10	1	VV	0.036	0.287
5	2.837	0.86	0	VV	0.032	0.118
6	2.885	0.68	0	VV	0.024	0.093
7	2.945	1.24	0	VV	0.052	0.170
8	3.169	5.53	1	VV	0.053	0.756
9	3.272	2.00	1	VV	0.044	0.273
10	3.344	4.42	1	VV	0.067	0.605
11	3.437	3.47	1	VV	0.044	0.474
12	3.486	11.09	1	VV	0.130	1.517
13	3.718	7.85	1	VV	0.107	1.074
14	3.850	2.34	1	VV	0.041	0.321
15	3.916	2.44	1	VV	0.054	0.333
16	4.020	2.87	2	VV	0.024	0.393
17	4.082	2.19	0	VB	0.081	0.300
18	4.239	2.57	1	BV	0.056	0.351
19	4.403	1.51	0	VV	0.049	0.206
20	4.444	1.36	0	VB	0.042	0.186
21	4.590	0.91	0	BV	0.061	0.125
22	4.701	1.77	0	VV	0.067	0.242
23	4.850	0.97	0	VV	0.033	0.132
24	4.911	0.53	0	VV	0.030	0.072
25	4.967	0.44	0	VV	0.028	0.061
26	5.069	1.34	1	VV	0.027	0.184
27	5.376	666.87	439	VV	0.025	91.184

Total area = 731.35

Area Percent Report

: W:\GC_45\GC 45 DATA\2014\141030\14103067.D Data File Name

Page Number

Operator : 682

: GC 45 Instrument

Report Created on: 04 Nov 14

Sample Name

: 14-10-2234-2

Vial Number

: Vial 67

Injection Number: 1

: 67 Sequence Line

Instrument Method: C:\CHEM32\1\METHODS\

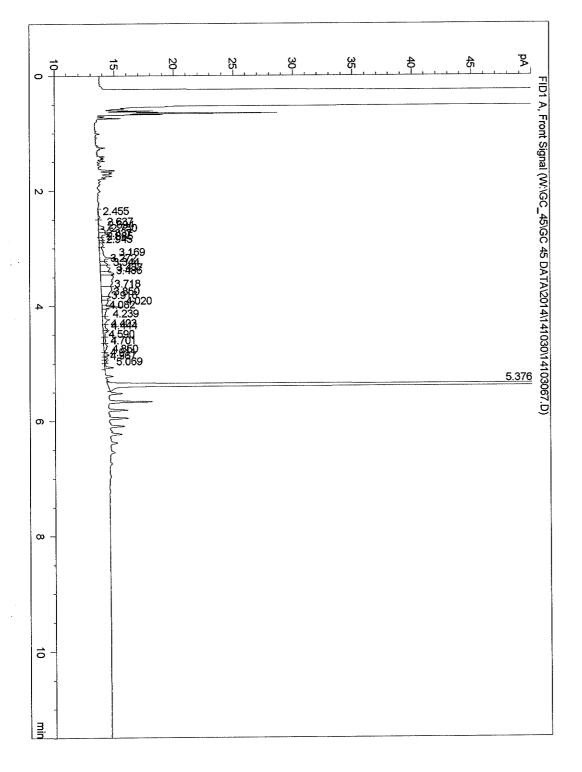
Acquired on

: 31 Oct 14

11:16 am 02:51 pm

Analysis Method : 8015B.MTH

Software Revision: Rev. B.03.02 [341] Copyright © Agilent Technologies



RAW DATA SHEET FOR METHOD: EPA 8015B (M)

WORK ORDER: INSTRUMENT:

14-10-2234

GC 45

EPA 3510C

EXTRACTION:

D/T EXTRACTED: 2014-10-30 00:00

ANALYZED BY:

D/T ANALYZED:

682 2014-10-31 11:35

REVIEWED BY: D/T REVIEWED:

DATA FILE:

W:\GC 45\GC 45 DATA\2014\141030\14103068.D\14103068

3

CLIENT SAMPLE NUMBER: ES118

LCS/MB BATCH:

141030B06 MS/MSD BATCH:

ug/L

SAMPLE VOLUME / WEIGHT:

DEFAULT: 500.00 ml / ACTUAL: 500.00 ml

FINAL VOLUME / WEIGHT:

DEFAULT: 5.00 ml / ACTUAL: 2.50 ml

ADJUSTMENT RATIO TO PF:

0.50

COMMENT:

UNITS:

Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, a

qualified with a "J" flag. TPH as Diesel is quantified in the carbon range C10-C28.

COMPOUND

INI. CONC DF

1.00

CONC 16.2

<u>DL</u>

LOQ LOD 25

QUAL

TPH as Diesel

3240

11

12

bJ

Area Percent Report

Data File Name : W:\GC_45\GC 45 DATA\2014\141030\14103068.D

Page Number : 1

Vial Number : Vial 68 Operator : 682

Operator : 682
Instrument : GC 45
Sample Name : 14-10-2234-3 Injection Number: 1 Sequence Line : 68

Instrument Method: C:\CHEM32\1\METHODS\ ->

Acquired on : 31 Oct 14 11:35 am

Analysis Method : 8015B.MTH Report Created on: 04 Nov 14 02:53 pm

Software Revision: Rev. B.03.02 [341] Copyright © Agilent Technologies

Sig. 1 in W:\GC 45\GC 45 DATA\2014\141030\ \rightarrow

	Ret Time	Area	Height			
•	-	0.55		VV	0.050	0.091
1 2	2.457 2.601	0.48		VV	0.030	0.080
3	2.640	0.40		VV	0.021	0.101
4	2.689	1.65		VV	0.024	0.276
5	2.752	1.54		VV	0.038	0.257
6	2.837	0.65		VV	0.034	0.108
7	2.885	0.64		VV	0.033	0.107
8	2.942	0.65		VV	0.043	0.109
9	3.101	1.29		vv	0.073	0.216
10	3.169	3.26		VV	0.026	0.545
11	3.342	3.49		VV	0.110	0.583
12	3.437	1.96		VV	0.049	0.327
13	3.481	1.65		VV	0.040	0.275
14	3.527	1.70		VV	0.042	0.284
15	3.596	3.74		VV	0.095	0.624
16	3.684	1.55	1	VV	0.047	0.259
17	3.755	3.47	1	VV	0.089	0.579
18	3.851	1.60	1	VV	0.045	0.268
19	3.916	0.61	0	VV	0.024	0.103
20	3.944	1.04	0	VV	0.047	0.174
21	4.020	2.38	2	VV	0.022	0.398
22	4.089	1.00	0	VV	0.054	0.167
23	4.132	0.32	0	VV	0.024	0.053
24	4.160	0.20	0	VV	0.022	0.033
25	4.239	2.07	1	VV	0.056	0.346
26	4.357	0.79	0	VV	0.044	0.132
27	4.406	1.51	0	VV	0.055	0.253
28	4.596	0.74		VV	0.065	0.124
29	4.723	1.18	0	VV	0.054	0.198
30	4.737	1.47		VV	0.030	0.246
31	4.851	0.96		VV	0.037	0.161
32	4.913	0.71		VV	0.030	0.118
33	4.972	0.41		VV	0.023	0.069
34	5.024	0.16		VV	0.023	0.027
35	5.073	0.91		VV	0.024	0.153
36	5.270	0.06		VM	0.010	0.010
37	5.379	551.80	370	VV	0.025	92.147

Area Percent Report

Data File Name : W:\GC_45\GC 45 DATA\2014\141030\14103068.D

Page Number : 2

Operator : 682 Vial Number : Vial 68

Instrument : GC 45 Injection Number : 1 Sample Name : 14-10-2234-3 Sequence Line : 68

Instrument Method: C:\CHEM32\1\METHODS\ ->

Acquired on : 31 Oct 14 11:35 am

Report Created on: 04 Nov 14 02:53 pm Analysis Method : 8015B.MTH

Software Revision: Rev. B.03.02 [341] Copyright © Agilent Technologies

Total area = 598.83

Area Percent Report

Data File Name : W:\GC 45\GC 45 DATA\2014\141030\14103068.D

Page Number :

Operator : 682 Vial Number : Vial 68

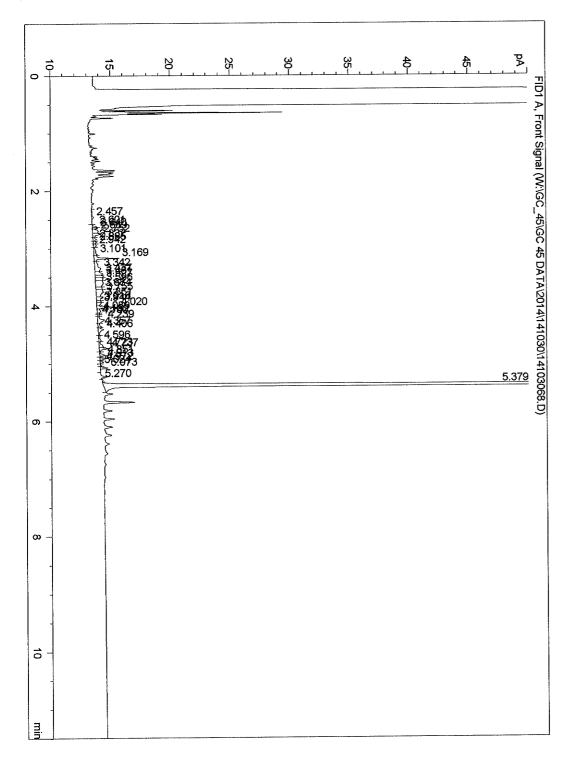
Instrument : GC 45 Injection Number : 1
Sample Name : 14-10-2234-3 Sequence Line : 68

Instrument Method: C:\CHEM32\1\METHODS\ ->

Acquired on : 31 Oct 14 11:35 am

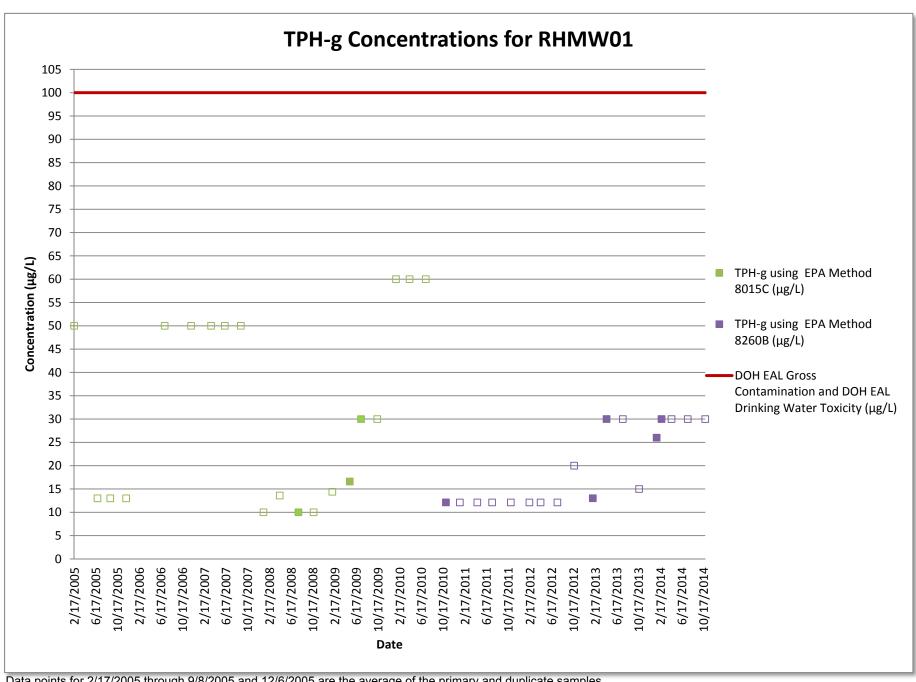
Report Created on: 04 Nov 14 02:53 pm Analysis Method : 8015B.MTH

Software Revision: Rev. B.03.02 [341] Copyright © Agilent Technologies



APPENDIX D Historical Groundwater Exceedances



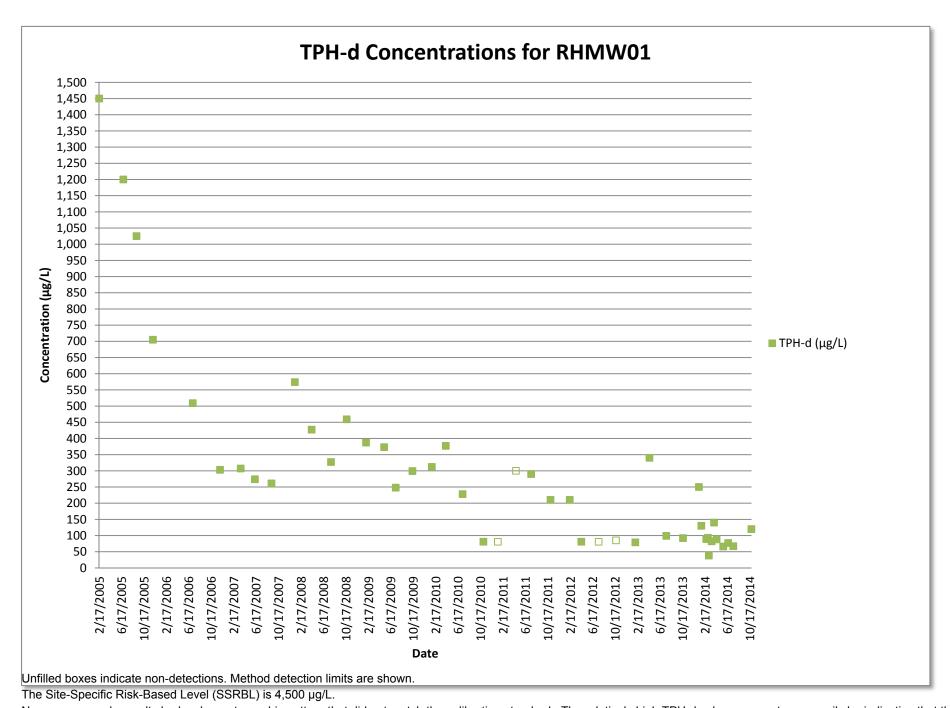


Data points for 2/17/2005 through 9/8/2005 and 12/6/2005 are the average of the primary and duplicate samples.

Possible laboratory contamination for 10/22/2012, 10/21/2013, and 1/28/2014 sampling events.

Unfilled boxes indicate non-detections. Method detection limits are shown.

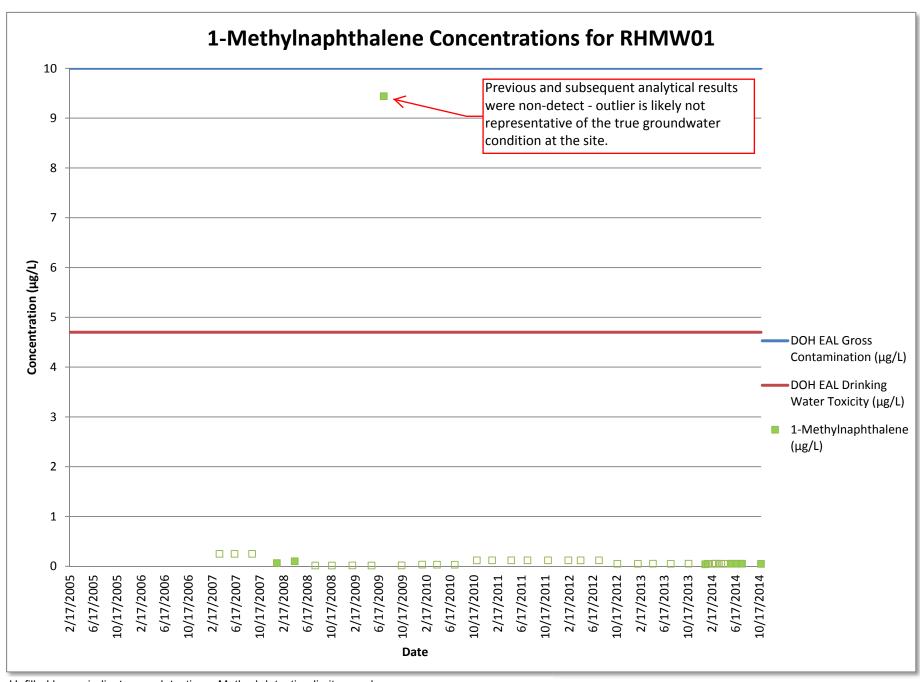




Numerous sample results had a chromatographic pattern that did not match the calibration standard. The relatively high TPH-d values may not necessarily be indicative that there is diesel fuel or other petroleum products in the well.

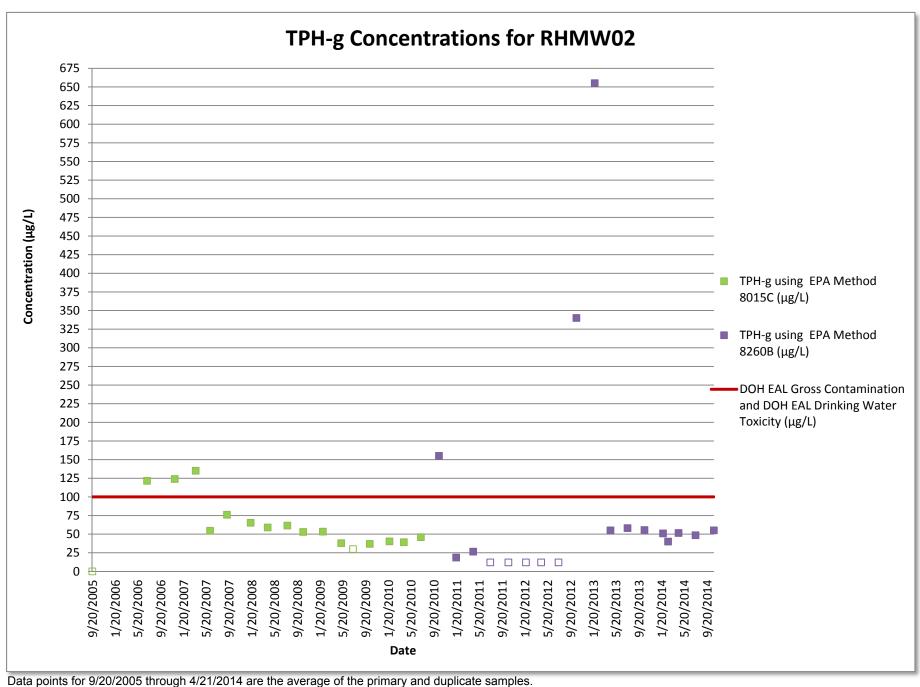
Data points for 2/17/2005 through 9/8/2005 and 12/6/2005 are the average of the primary and duplicate samples.





Unfilled boxes indicate non-detections. Method detection limits are shown.



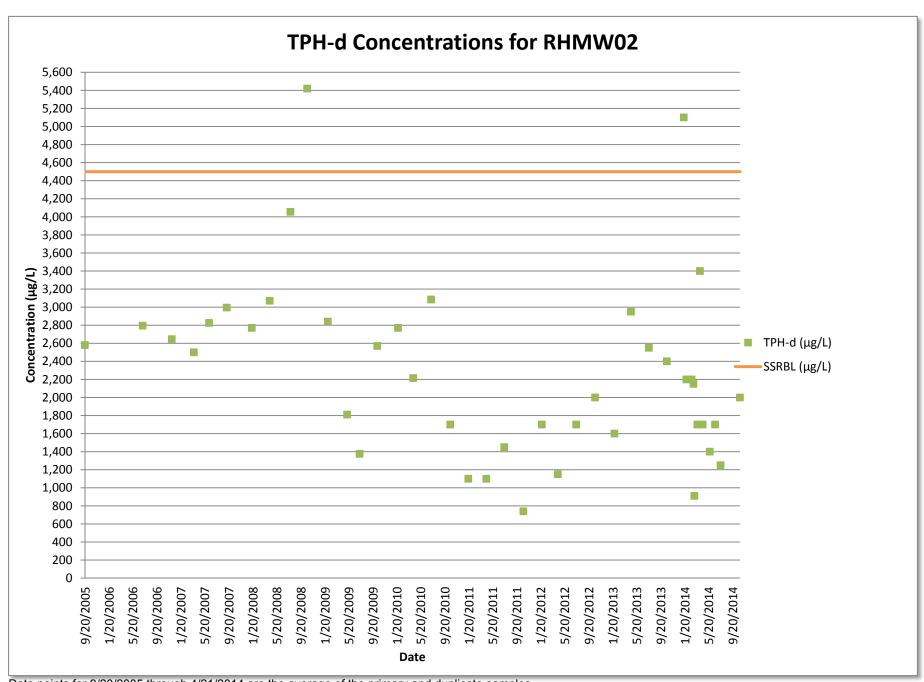


Possible laboratory contamination for 10/21/2013 and 1/28/2014 sampling events.

Unfilled boxes indicate non-detections. Method detection limits are shown.

Primary sample results are shown for 1/26/2012 and 7/18/2012; all other concentrations are the average of the primary and duplicate sample results.

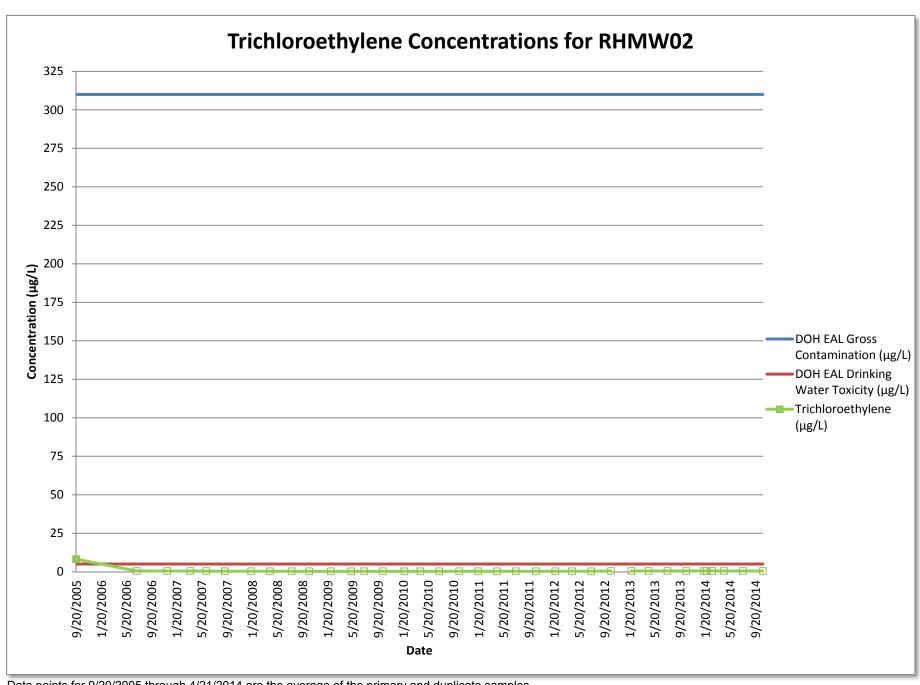




Data points for 9/20/2005 through 4/21/2014 are the average of the primary and duplicate samples.

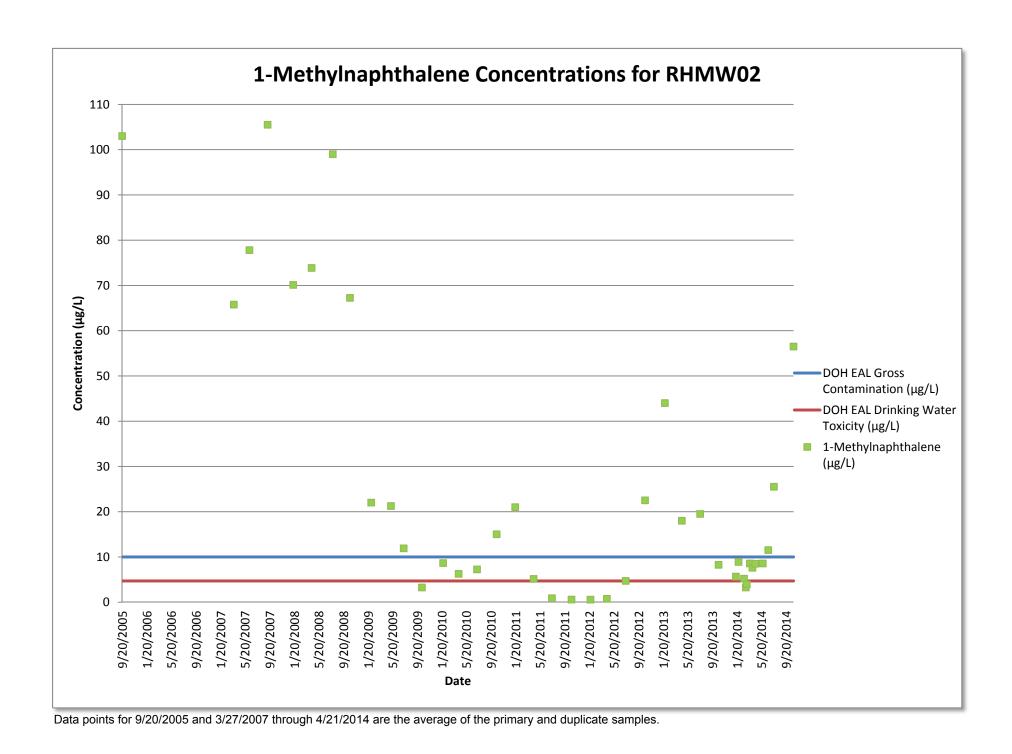
Numerous sample results had a chromatographic pattern that did not match the calibration standard. The relatively high TPH-d values may not necessarily be indicative that there is diesel fuel or other petroleum products in the well.



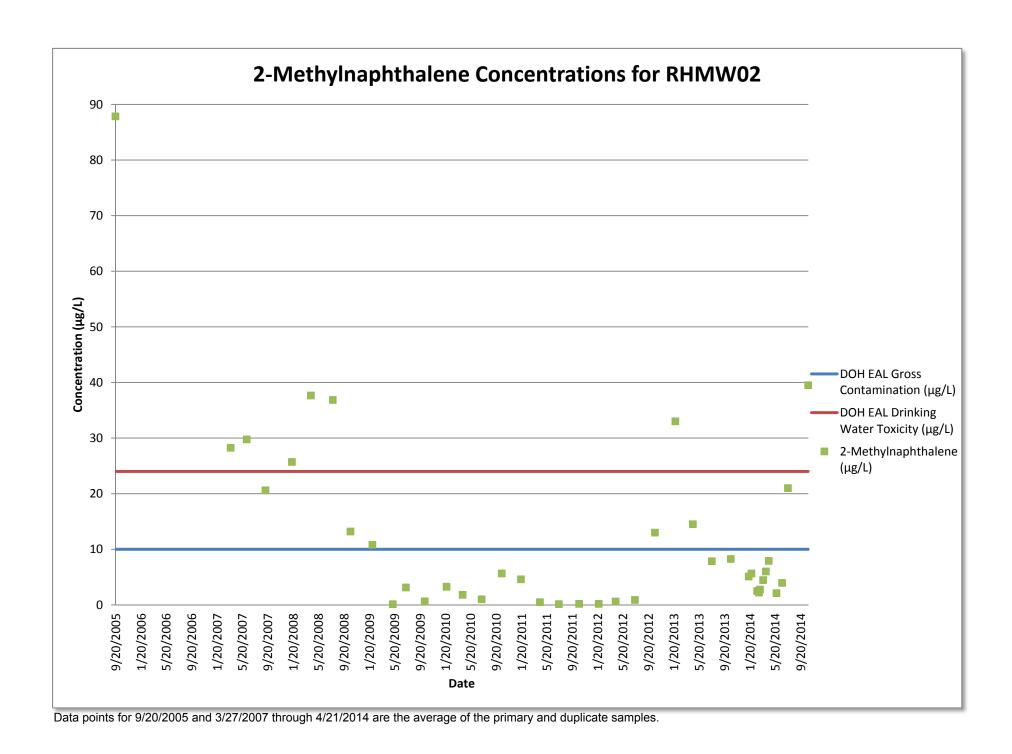


Data points for 9/20/2005 through 4/21/2014 are the average of the primary and duplicate samples. Unfilled boxes indicate non-detections. Method detection limits are shown.

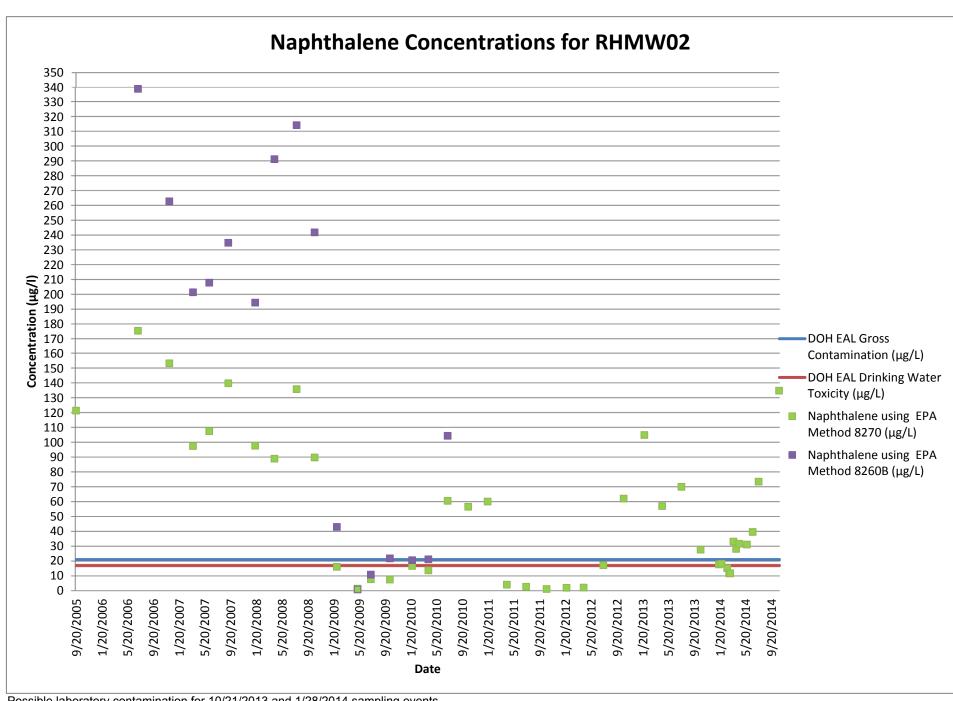






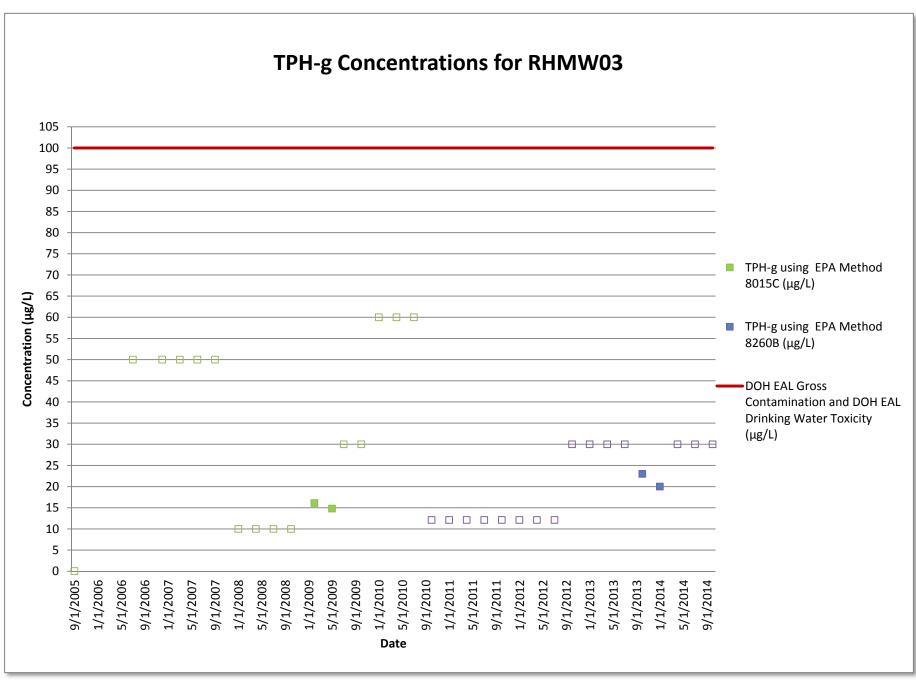






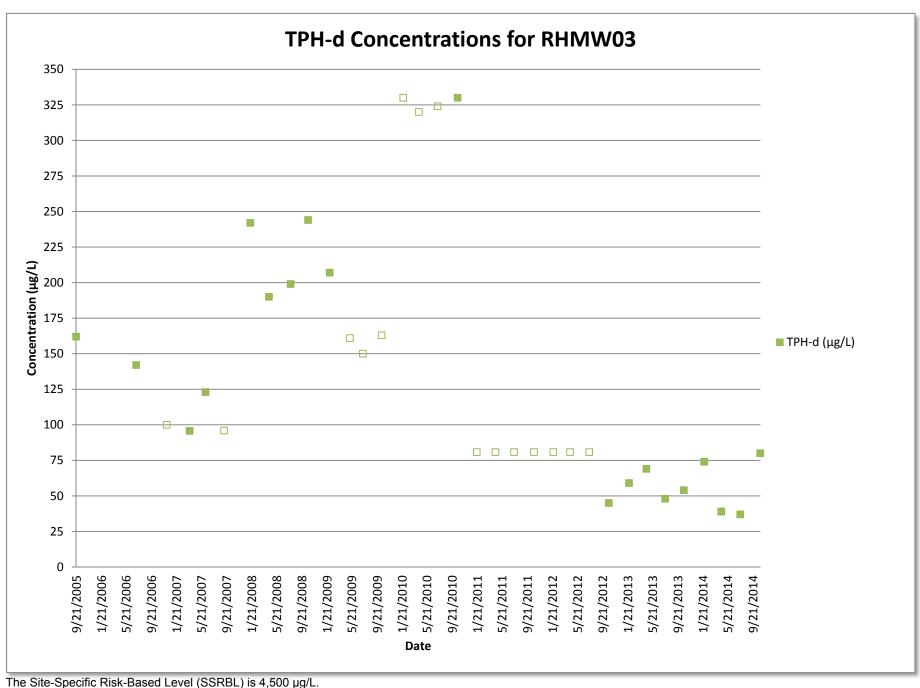
Possible laboratory contamination for 10/21/2013 and 1/28/2014 sampling events. Unfilled boxes indicate non-detections. Method detection limits are shown.





Possible laboratory contamination for 10/21/2013 and 1/28/2014 sampling events. Unfilled boxes indicate non-detections. Method detection limits are shown.

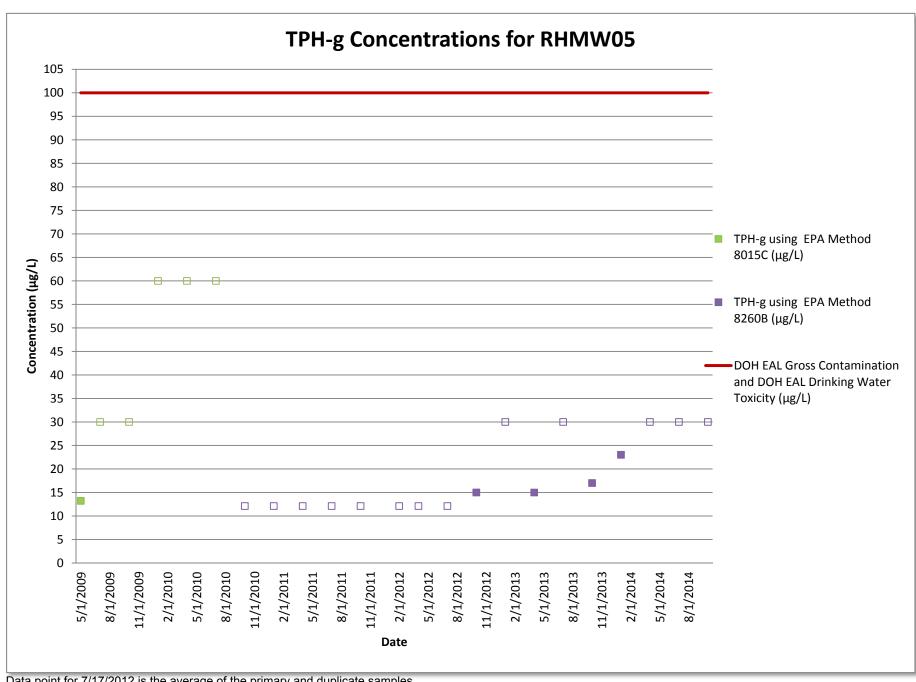




Unfilled boxes indicate non-detections. Method detection limits are shown.

Numerous sample results had a chromatographic pattern that didn't match the calibration standard. The relatively high TPH-d values may not necessarily be indicative that there is diesel fuel or other petroleum products in the well.



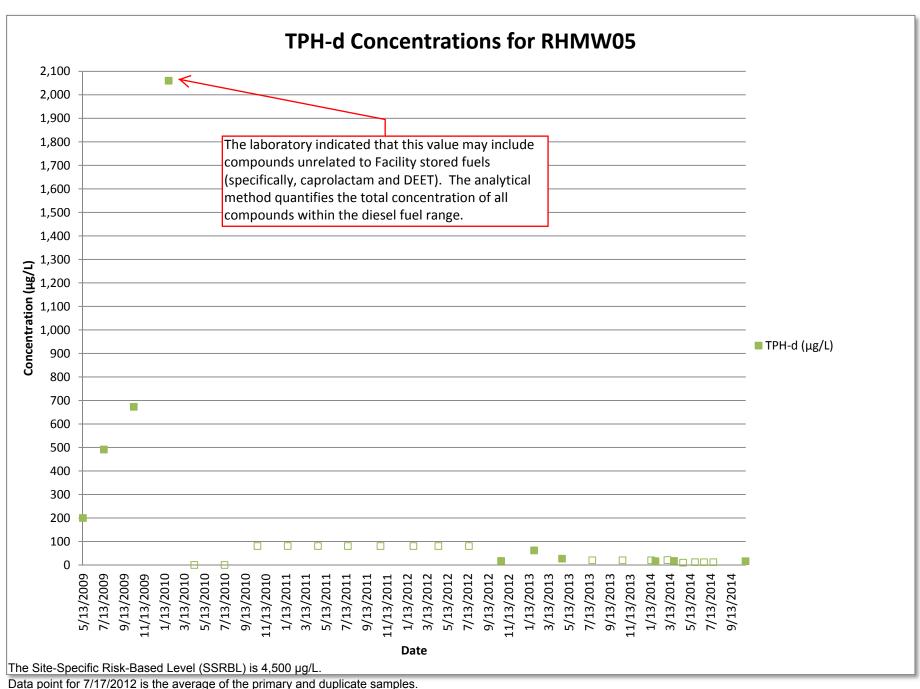


Data point for 7/17/2012 is the average of the primary and duplicate samples.

Unfilled boxes indicate non-detections. Method detection limits are shown.

Possible laboratory contamination for 10/22/2012, 10/22/2013, and 1/29/2014 sampling events.



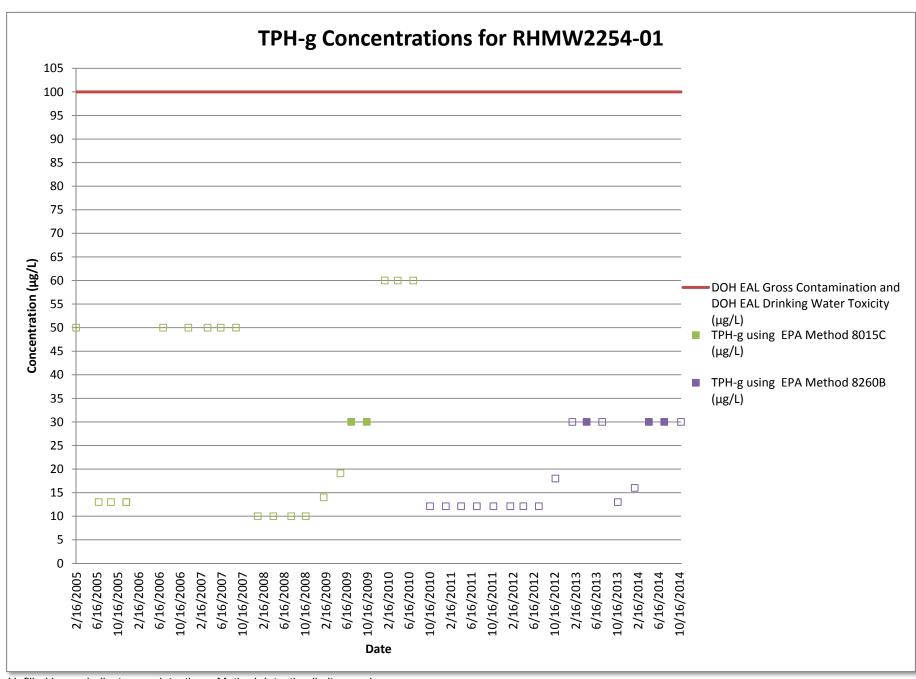


Data point for 7/17/2012 is the average of the primary and duplicate samples.

Unfilled boxes indicate non-detections. Method detection limits are shown.

Numerous sample results had a chromatographic pattern that did not match the calibration standard. The relatively high TPH-d values may not necessarily be indicative that there is diesel fuel or other petroleum products in the well.

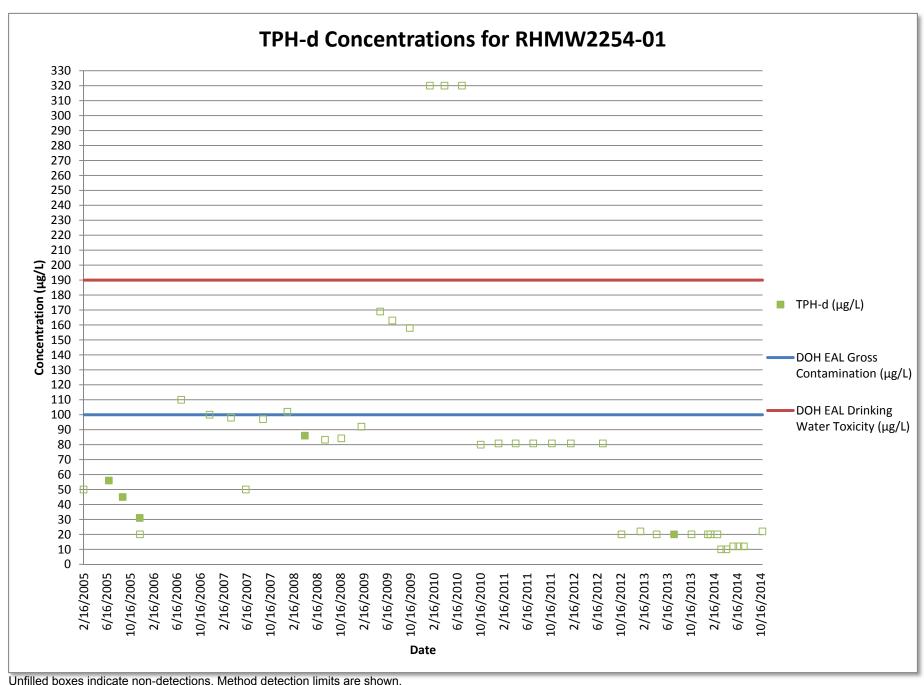




Unfilled boxes indicate non-detections. Method detection limits are shown.

Possible laboratory contamination for 10/22/2012, 10/22/2013, and 1/29/2014 sampling events.





Laboratory data rejected for 1/15/2008 sampling event.

Numerous sample results had a chromatographic pattern that did not match the calibration standard. The relatively high TPH-d values may not necessarily be indicative that there is diesel fuel or other petroleum products in the well.



APPENDIX E

Cumulative Groundwater Results (included on attached CD)

